

# Robust Recovery of Signals From a Structured Union of Subspaces

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**Abstract**—Traditional sampling theories consider the problem of reconstructing an unknown signal  $x$  from a series of samples. A prevalent assumption which often guarantees recovery from the given measurements is that  $x$  lies in a known subspace. Recently, there has been growing interest in nonlinear but structured signal models, in which  $x$  lies in a union of subspaces. In this paper, we develop a general framework for robust and efficient recovery of such signals from a given set of samples. More specifically, we treat the case in which  $x$  lies in a sum of  $k$  subspaces, chosen from a larger set of  $m$  possibilities. The samples are modeled as inner products with an arbitrary set of sampling functions. To derive an efficient and robust recovery algorithm, we show that our problem can be formulated as that of recovering a block-sparse vector whose nonzero elements appear in fixed blocks. We then propose a mixed  $\ell_2/\ell_1$  program for block sparse recovery. Our main result is an equivalence condition under which the proposed convex algorithm is guaranteed to recover the original signal. This result relies on the notion of block restricted isometry property (RIP), which is a generalization of the standard RIP used extensively in the context of compressed sensing. Based on RIP, we also prove stability of our approach in the presence of noise and modeling errors. A special case of our framework is that of recovering multiple measurement vectors (MMV) that share a joint sparsity pattern. Adapting our results to this context leads to new MMV recovery methods as well as equivalence conditions under which the entire set can be determined efficiently.

**Index Terms**—Block restricted isometry property, block sparsity, compressed sensing, mixed-norm recovery, multiple measurement vectors (MMV), union of linear subspaces.

## I. INTRODUCTION

SAMPLING theory has a rich history dating back to Cauchy. Undoubtedly, the sampling theorem that had the most impact on signal processing and communications is that associated with Whittaker, Kotelníkov, and Shannon [1], [2]. Their famous result is that a bandlimited function  $x(t)$  can be recovered from its uniform samples as long as the sampling rate exceeds the Nyquist rate, corresponding to twice the highest frequency of the signal [3]. More recently, this basic theorem has been extended to include more general classes of signal spaces. In par-

Manuscript received July 25, 2008; revised March 27, 2009. Current version published October 21, 2009. This work was supported in part by the Israel Science Foundation under Grant 1081/07 and by the European Commission in the framework of the FP7 Network of Excellence in Wireless COMMunications NEWCOM++ under Contract 216715. The material in this paper was presented in part at the 16th International Conference on Digital Signal Processing (DSP'2009), Santorini, Greece, July 2009.

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Communicated by J. Romberg, Associate Editor for Signal Processing.

Color versions of Figures 1–5 in this paper are available online at <http://ieeexplore.ieee.org>.

Digital Object Identifier 10.1109/TIT.2009.2030471

ticular, it can be shown that under mild technical conditions, a signal  $x$  lying in a given subspace can be recovered exactly from its linear generalized samples using a series of filtering operations [4]–[7].

Recently, there has been growing interest in nonlinear signal models in which the unknown  $x$  does not necessarily lie in a subspace. In order to ensure recovery from the samples, some underlying structure is needed. A general model that captures many interesting cases is that in which  $x$  lies in a union of subspaces. In this setting,  $x$  resides in one of a set of given subspaces  $\mathcal{V}_i$ , however, *a priori* it is not known in which one. A special case of this framework is the problem underlying the field of compressed sensing (CS), in which the goal is to recover a length- $N$  vector  $\mathbf{x}$  from  $n < N$  linear measurements, where  $\mathbf{x}$  has no more than  $k$  nonzero elements in some basis [8], [9]. Many algorithms have been proposed in the literature in order to recover  $\mathbf{x}$  in a stable and efficient manner [9]–[13]. A variety of conditions have been developed to ensure that these methods recover  $\mathbf{x}$  exactly. One of the main tools in this context is the restricted isometry property (RIP) [9], [14], [15]. In particular, it can be shown that if the measurement matrix satisfies the RIP with an appropriate constant, then  $\mathbf{x}$  can be recovered by solving an  $\ell_1$  minimization algorithm.

Another special case of a union of subspaces is the setting in which the unknown signal  $x = x(t)$  has a multiband structure, so that its Fourier transform consists of a limited number of bands at unknown locations [16], [17]. By formulating this problem within the framework of CS, explicit sub-Nyquist sampling and reconstruction schemes were developed in [16], [17] that ensure perfect recovery at the minimal possible rate. This setup was recently generalized in [18], [19] to deal with sampling and reconstruction of signals that lie in a finite union of shift-invariant subspaces. By combining ideas from standard sampling theory with CS results [20], explicit low-rate sampling and recovery methods were developed for such signal sets. In [21], an extension was considered to a special case of an infinite union of shift-invariant subspaces. The infinite union is a result of the fact that each generator of the space has an unknown time delay. Another example of a union of subspaces is the set of finite rate of innovation signals [21]–[23], that are modeled as a weighted sum of shifts of a given generating function, where the shifts are unknown.

In this paper, our goal is to develop a unified framework for efficient recovery of signals that lie in a structured union of subspaces. Our emphasis is on computationally efficient methods that are stable in the presence of noise and modeling errors. In contrast to our previous work [16]–[19], [21], here we consider unions of finite-dimensional subspaces. Specifically, we restrict

our attention to the case in which  $x$  resides in a sum of  $k$  subspaces, chosen from a given set of  $m$  subspaces  $\mathcal{A}_j$ ,  $1 \leq j \leq m$ . However, which subspaces comprise the sum is unknown. This setting is a special case of the more general union model treated in [24], [25]. Conditions under which unique and stable sampling are possible were developed in [24], [25]. However, no concrete algorithm was provided to recover such a signal from a given set of samples in a stable and efficient manner. Here, we propose a convex optimization algorithm that will often recover the true underlying  $x$ , and develop explicit conditions under which perfect recovery is guaranteed. Furthermore, we prove that our method is stable and robust in the sense that the reconstruction error is bounded in the presence of noise and mismatching; namely, when  $x$  does not lie exactly in the union. Our results rely on a generalization of the RIP which fits the union setting we treat here.

Our first contribution is showing that the problem of recovering  $x$  in a structured union of subspaces can be cast as a sparse recovery problem, in which it is desired to recover a sparse vector  $\mathbf{c}$  that has a particular sparsity pattern: the nonzero values appear in fixed blocks. We refer to such a model as block sparsity. Clearly, any block-sparse vector is also sparse in the standard sense. However, by exploiting the block structure of the sparsity pattern, recovery may be possible under more general conditions.

Next, we develop a concrete algorithm to determine a block-sparse vector from given measurements, which is based on minimizing a mixed  $\ell_2/\ell_1$  norm. This problem can be cast as a convex second-order cone program (SOCP), and solved efficiently using standard software packages. A mixed norm approach for block-sparse recovery was also considered in [26], [27]. By analyzing the measurement operator's null space, it was shown that asymptotically, as the signal length grows to infinity, and under ideal conditions (no noise or modeling errors), perfect recovery is possible with high probability. However, no robust equivalence results were established between the output of the algorithm and the true block-sparse vector for a given finite-length measurement vector, or in the presence of noise and mismatching.

Generalizing the concept of RIP to our setting, we introduce the block RIP, which is a less stringent requirement. We then prove that if the measurement matrix satisfies the block RIP, then our proposed convex algorithm will recover the underlying block sparse signal. Furthermore, under block RIP, our algorithm is stable in the presence of noise and mismatching errors. Using ideas similar to [13], [28] we then prove that random matrices satisfy the block RIP with overwhelming probability. Moreover, the probability to satisfy the block RIP is substantially larger than that of satisfying the standard RIP. These results establish that a signal  $x$  that lies in a finite structured union can be recovered efficiently and stably with overwhelming probability if the measurement matrix is constructed from a random ensemble.

An interesting special case of the block-sparse model is the multiple measurement vector (MMV) problem, in which there is a set of unknown vectors that share a joint sparsity pattern. MMV recovery algorithms were studied in [20], [29]–[32]. Equivalence results based on mutual coherence for a mixed

$\ell_p/\ell_1$  program were derived in [30]. These results turn out to be the same as that obtained from a single measurement problem. This is in contrast to the fact that in practice, MMV methods tend to outperform algorithms that treat each of the vectors separately. In order to develop meaningful equivalence results, we cast the MMV problem as one of block-sparse recovery. Our mixed  $\ell_2/\ell_1$  method translates into minimizing the sum of the  $\ell_2$  row-norms of the unknown matrix representing the MMV set. Our general results lead to RIP-based equivalence conditions for this algorithm. Furthermore, our framework suggests a different type of sampling method for MMV problems which tends to increase the recovery rate. The equivalence condition we obtain in this case is stronger than the single measurement setting. As we show, this method leads to superior recovery rate when compared with other popular MMV algorithms.

The remainder of the paper is organized as follows. In Section II, we describe the general problem of sampling from a union of subspaces. The relationship between our problem and that of block-sparse recovery is developed in Section III. In Section IV, we explore stability and uniqueness issues which leads to the definition of block RIP. We also present a nonconvex optimization algorithm with combinatorial complexity whose solution is the true unknown  $x$ . A convex relaxation of this algorithm is proposed in Section V. We then derive equivalence conditions based on block RIP. The concept of block RIP is further used to establish robustness and stability of our algorithm in the presence of noise and modeling errors. This approach is specialized to MMV sampling in Section VI. Finally, in Section VII, we prove that random ensembles tend to satisfy the block RIP with high probability.

Throughout the paper, we denote vectors in an arbitrary Hilbert space  $\mathcal{H}$  by lower case letters e.g.,  $x$ , and sets of vectors in  $\mathcal{H}$  by calligraphic letters, e.g.,  $\mathcal{S}$ . Vectors in  $\mathbb{R}^N$  are written as boldface lowercase letters e.g.,  $\mathbf{x}$ , and matrices as boldface uppercase letters e.g.,  $\mathbf{A}$ . The identity matrix of appropriate dimension is written as  $\mathbf{I}$  or  $\mathbf{I}_d$  when the dimension is not clear from the context, and  $\mathbf{A}^T$  is the transpose of the matrix  $\mathbf{A}$ . The  $i$ th element of a vector  $\mathbf{x}$  is denoted by  $\mathbf{x}(i)$ . Linear transformations from  $\mathbb{R}^n$  to  $\mathcal{H}$  are written as upper case letters  $A : \mathbb{R}^n \rightarrow \mathcal{H}$ . The adjoint of  $A$  is written as  $A^*$ . The standard Euclidean norm is denoted  $\|\mathbf{x}\|_2 = \sqrt{\mathbf{x}^T \mathbf{x}}$  and  $\|\mathbf{x}\|_1 = \sum_i |\mathbf{x}(i)|$  is the  $\ell_1$  norm of  $\mathbf{x}$ . The Kronecker product between matrices  $\mathbf{A}$  and  $\mathbf{B}$  is denoted  $\mathbf{A} \otimes \mathbf{B}$ , and  $\text{vec}(\mathbf{Y})$  is the vector obtained by concatenating the columns of  $\mathbf{Y}$ . The following variables are used in the sequel:  $n$  is the number of samples,  $N$  is the length of the input signal  $\mathbf{x}$  when it is a vector,  $k$  is the sparsity or block sparsity (to be defined later on) of a vector  $\mathbf{c}$ , and  $m$  is the number of subspaces. For ease of notation, we assume throughout that all scalars are defined over the field of real numbers; however, the results are also valid over the complex domain with appropriate modifications.

## II. UNION OF SUBSPACES

### A. Subspace Sampling

Traditional sampling theory deals with the problem of recovering an unknown signal  $x \in \mathcal{H}$  from a set of  $n$  samples  $y_i = f_i(x)$  where  $f_i(x)$  is some function of  $x$ . The signal  $x$  can

be a function of time  $x = x(t)$ , or can represent a finite-length vector  $x = \mathbf{x}$ . The most common type of sampling is linear sampling in which

$$y_i = \langle s_i, x \rangle, \quad 1 \leq i \leq n \quad (1)$$

for a set of functions  $s_i \in \mathcal{H}$  [4], [33]–[39]. Here  $\langle x, y \rangle$  denotes the standard inner product on  $\mathcal{H}$ . For example, if  $\mathcal{H} = L_2$  is the space of real finite-energy signals then

$$\langle x, y \rangle = \int_{-\infty}^{\infty} x(t)y(t)dt. \quad (2)$$

When  $\mathcal{H} = \mathbb{R}^N$  for some  $N$

$$\langle \mathbf{x}, \mathbf{y} \rangle = \sum_{i=1}^N \mathbf{x}(i)\mathbf{y}(i). \quad (3)$$

Nonlinear sampling is treated in [40]. However, here our focus will be on the linear case.

When  $\mathcal{H} = \mathbb{R}^N$  the unknown  $x = \mathbf{x}$  as well as the sampling functions  $s_i = \mathbf{s}_i$  are vectors in  $\mathbb{R}^N$ . Therefore, the samples can be written conveniently in matrix form as  $\mathbf{y} = \mathbf{S}^T \mathbf{x}$ , where  $\mathbf{S}$  is the matrix with columns  $\mathbf{s}_i$ . In the more general case in which  $\mathcal{H} = L_2$  or any other abstract Hilbert space, we can use the set transformation notation in order to conveniently represent the samples. A set transformation  $S : \mathbb{R}^n \rightarrow \mathcal{H}$  corresponding to sampling vectors  $\{s_i \in \mathcal{H}, 1 \leq i \leq n\}$  is defined by

$$\mathbf{S}\mathbf{c} = \sum_{i=1}^n \mathbf{c}(i)s_i \quad (4)$$

for all  $\mathbf{c} \in \mathbb{R}^n$ . From the definition of the adjoint, if  $\mathbf{c} = S^*x$ , then  $\mathbf{c}(i) = \langle s_i, x \rangle$ . Note that when  $\mathcal{H} = \mathbb{R}^N$ ,  $S = \mathbf{S}$  and  $S^* = \mathbf{S}^T$ . Using this notation, we can always express the samples as

$$\mathbf{y} = S^*x \quad (5)$$

where  $S$  is a set transformation for arbitrary  $\mathcal{H}$ , and an appropriate matrix when  $\mathcal{H} = \mathbb{R}^N$ .

Our goal is to recover  $x$  from the samples  $\mathbf{y} \in \mathbb{R}^n$ . If the vectors  $s_i$  do not span the entire space  $\mathcal{H}$ , then there are many possible signals  $x$  consistent with  $\mathbf{y}$ . More specifically, if we define by  $\mathcal{S}$  the sampling space spanned by the vectors  $s_i$ , then clearly  $S^*v = 0$  for any  $v \in \mathcal{S}^\perp$ . Therefore, if  $\mathcal{S}^\perp$  is not the trivial space then adding such a vector  $v$  to any solution  $x$  of (5) will result in the same samples  $\mathbf{y}$ . However, by exploiting prior knowledge on  $x$ , in many cases uniqueness can be guaranteed. A prior very often assumed is that  $x$  lies in a given subspace  $\mathcal{A}$  of  $\mathcal{H}$  [4]–[7]. If  $\mathcal{A}$  and  $\mathcal{S}$  have the same finite dimension, and  $\mathcal{S}^\perp$  and  $\mathcal{A}$  intersect only at the 0 vector, then  $x$  can be perfectly recovered from the samples  $\mathbf{y}$  [6], [7], [41].

### B. Union of Subspaces

When subspace information is available, perfect reconstruction can often be guaranteed. Furthermore, recovery can be implemented by a simple linear transformation of the given samples (5). However, there are many practical scenarios in which we are given prior information about  $x$  that is not necessarily in the form of a subspace. One such case studied in detail in [41]

is that in which  $x$  is known to be smooth. Here we focus our attention on the setting where  $x$  lies in a union of subspaces

$$\mathcal{U} = \bigcup_i \mathcal{V}_i \quad (6)$$

where each  $\mathcal{V}_i$  is a subspace. Thus,  $x$  belongs to one of the  $\mathcal{V}_i$ , but we do not know *a priori* to which one [24], [25]. Note that the set  $\mathcal{U}$  is no longer a subspace. Indeed, if  $\mathcal{V}_i$  is, for example, a one-dimensional space spanned by the vector  $\mathbf{v}_i$ , then  $\mathcal{U}$  contains vectors of the form  $\alpha\mathbf{v}_i$  for some  $i$  but does not include their linear combinations. Our goal is to recover a vector  $x$  lying in a union of subspaces, from a given set of samples. In principle, if we knew which subspace  $x$  belonged to, then reconstruction can be obtained using standard sampling results. However, here the problem is more involved because conceptually we first need to identify the correct subspace and only then can we recover the signal within the space.

Previous work on sampling over a union focused on invertibility and stability results [24], [25]. In contrast, here, our main interest is in developing concrete recovery algorithms that are provably robust. To achieve this goal, we limit our attention to a subclass of (6) for which stable recovery algorithms can be developed and analyzed. Specifically, we treat the case in which each  $\mathcal{V}_i$  has the additional structure

$$\mathcal{V}_i = \bigoplus_{|j|=k} \mathcal{A}_j \quad (7)$$

where  $\{\mathcal{A}_j, 1 \leq j \leq m\}$  are a given set of disjoint subspaces, and  $|j| = k$  denotes a sum over  $k$  indices. Thus, each subspace  $\mathcal{V}_i$  corresponds to a different choice of  $k$  subspaces  $\mathcal{A}_j$  that comprise the sum. We assume throughout the paper that  $m$  and the dimensions  $d_i = \dim(\mathcal{A}_i)$  of the subspaces  $\mathcal{A}_i$  are finite. Given  $n$  samples

$$\mathbf{y} = S^*x \quad (8)$$

and the knowledge that  $x$  lies in exactly one of the subspaces  $\mathcal{V}_i$ , we would like to recover the unknown signal  $x$ . In this setting, there are  $\binom{m}{k}$  possible subspaces comprising the union.

An alternative interpretation of our model is as follows. Given an observation vector  $\mathbf{y}$ , we seek a signal  $x$  for which  $\mathbf{y} = S^*x$  and in addition  $x$  can be written as

$$x = \sum_{i=1}^k x_i \quad (9)$$

where each  $x_i$  lies in  $\mathcal{A}_j$  for some index  $j$ .

A special case is the standard CS problem in which  $x = \mathbf{x}$  is a vector of length  $N$ , that has a sparse representation in a given basis defined by an invertible matrix  $\mathbf{W}$ . Thus,  $\mathbf{x} = \mathbf{W}\mathbf{c}$  where  $\mathbf{c}$  is a sparse vector that has at most  $k$  nonzero elements. This fits our framework by choosing  $\mathcal{A}_i$  as the space spanned by the  $i$ th column of  $\mathbf{W}$ . In this setting  $m = N$ , and there are  $\binom{N}{k}$  subspaces comprising the union.

Another example is the block sparsity model [26], [42] in which  $\mathbf{x}$  is divided into equal-length blocks of size  $d$ , and at most  $k$  blocks can be nonzero. Such a vector can be described

in our setting with  $\mathcal{H} = \mathbb{R}^N$  by choosing  $\mathcal{A}_i$  to be the space spanned by the corresponding  $d$  columns of the identity matrix. Here  $m = N/d$  and there are  $\binom{N/d}{k}$  subspaces in the union.

A final example is the MMV problem [20], [29]–[32] in which our goal is to recover a matrix  $\mathbf{X}$  from measurements  $\mathbf{Y} = \mathbf{M}\mathbf{X}$ , for a given sampling matrix  $\mathbf{M}$ . The matrix  $\mathbf{X}$  is assumed to have at most  $k$  nonzero rows. Thus, not only is each column  $\mathbf{x}_i$   $k$ -sparse, but in addition the nonzero elements of  $\mathbf{x}_i$  share a joint sparsity pattern. This problem can be transformed into that of recovering a  $k$ -block sparse signal by stacking the rows of  $\mathbf{X}$  and  $\mathbf{Y}$ , leading to the relationship

$$\text{vec}(\mathbf{Y}^T) = (\mathbf{M} \otimes \mathbf{I})\text{vec}(\mathbf{X}^T). \quad (10)$$

The structure of  $\mathbf{X}$  leads to a vector  $\text{vec}(\mathbf{X}^T)$  that is  $k$ -block sparse.

### C. Problem Formulation and Main Results

Given  $k$  and the subspaces  $\mathcal{A}_i$ , we would like to address the following questions.

- 1) What are the conditions on the sampling vectors  $s_i, 1 \leq i \leq n$  in order to guarantee that the sampling is invertible and stable?
- 2) How can we recover the unique  $x$  (regardless of computational complexity)?
- 3) How can we recover the unique  $x$  in an efficient and stable manner?

The first question was addressed in [24], [25] in the more general context of unions of spaces (without requiring a particular structure such as (7)). However, no concrete methods were proposed in order to recover  $x$ . Here we provide efficient convex algorithms that recover  $x$  in a stable way for arbitrary  $k$  under appropriate conditions on the sampling functions  $s_i$  and the spaces  $\mathcal{A}_i$ .

Our results are based on an equivalence between the union of subspaces problem assuming (7) and that of recovering block-sparse vectors. This allows us to determine  $x$  from the given samples by first treating the problem of recovering a block  $k$ -sparse vector  $\mathbf{c}$  from a given set of measurements. This relationship is established in Section III. In the remainder of the paper we therefore focus on the block  $k$ -sparse model and develop our results in that context. In particular, we introduce a block RIP condition that ensures uniqueness and stability of our sampling problem. We then suggest an efficient convex optimization problem which approximates an unknown block-sparse vector  $\mathbf{c}$ . Based on block RIP we prove that  $\mathbf{c}$  can be recovered exactly in a stable way using the proposed optimization program. Furthermore, in the presence of noise and modeling errors, our algorithm can approximate the best block- $k$  sparse solution.

### III. CONNECTION WITH BLOCK SPARSITY

Consider the model of a signal  $x$  in the union of  $k$  out of  $m$  subspaces  $\mathcal{A}_i$ , with  $d_i = \dim(\mathcal{A}_i)$  as in (6) and (7). To write  $x$  explicitly, we choose a basis for each  $\mathcal{A}_i$ . Denoting by  $A_i : \mathbb{R}^{d_i} \rightarrow \mathcal{H}$  the set transformation corresponding to a basis for  $\mathcal{A}_i$ , any such  $x$  can be written as

$$x = \sum_{|i|=k} A_i \mathbf{c}_i \quad (11)$$

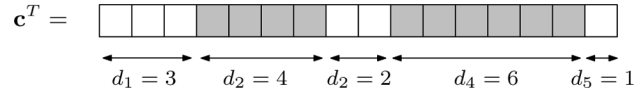


Fig. 1. A block-sparse vector  $\mathbf{c}$  over  $\mathcal{I} = \{d_1, \dots, d_5\}$ . The gray areas represent 10 nonzero entries which occupy two blocks.

where  $\mathbf{c}_i \in \mathbb{R}^{d_i}$  are the representation coefficients in  $\mathcal{A}_i$ , and  $|i| = k$  denotes a sum over a set of  $k$  indices. The choice of indices depend on the signal  $x$  and are unknown in advance.

To develop the equivalence with block sparsity, it is useful to introduce some further notation. First, we define  $A : \mathbb{R}^N \rightarrow \mathcal{H}$  as the set transformation that is a result of concatenating the different  $A_i$ , with

$$N = \sum_{i=1}^m d_i. \quad (12)$$

Next, we define the  $i$ th subblock  $\mathbf{c}[i]$  of a length- $N$  vector  $\mathbf{c}$  over  $\mathcal{I} = \{d_1, \dots, d_m\}$ . The  $i$ th subblock is of length  $d_i$ , and the blocks are formed sequentially so that

$$\mathbf{c}^T = [\underbrace{c_1 \dots c_{d_1}}_{\mathbf{c}[1]} \dots \underbrace{c_{N-d_m+1} \dots c_N}_{\mathbf{c}[m]}]^T. \quad (13)$$

We can then define  $A$  by

$$A\mathbf{c} = \sum_{i=1}^m A_i \mathbf{c}[i]. \quad (14)$$

When  $\mathcal{H} = \mathbb{R}^N$  for some  $N$ ,  $A_i = \mathbf{A}_i$  is a matrix and  $A = \mathbf{A}$  is the matrix obtained by column-wise concatenating  $\mathbf{A}_i$ . If for a given  $x$  the  $j$ th subspace  $\mathcal{A}_j$  does not appear in the sum (7), or equivalently in (11), then  $\mathbf{c}[j] = \mathbf{0}$ .

Any  $x$  in the union (6), (7) can be represented in terms of  $k$  of the bases  $A_i$ . Therefore, we can write  $x = A\mathbf{c}$  where there are at most  $k$  nonzero blocks  $\mathbf{c}[i]$ . Consequently, our union model is equivalent to the model in which  $x$  is represented by a sparse vector  $\mathbf{c}$  in an appropriate basis. However, the sparsity pattern here has a unique form which we will exploit in our conditions and algorithms: the nonzero elements appear in blocks.

*Definition 1:* A vector  $\mathbf{c} \in \mathbb{R}^N$  is called block  $k$ -sparse over  $\mathcal{I} = \{d_1, \dots, d_m\}$  if  $\mathbf{c}[i]$  is nonzero for at most  $k$  indices  $i$  where  $N = \sum_i d_i$ .

An example of a block-sparse vector with  $k = 2$  is depicted in Fig. 1. When  $d_i = 1$  for each  $i$ , block sparsity reduces to the conventional definition of a sparse vector. Denoting

$$\|\mathbf{c}\|_{0,\mathcal{I}} = \sum_{i=1}^m I(\|\mathbf{c}[i]\|_2 > 0) \quad (15)$$

where  $I(\|\mathbf{c}[i]\|_2 > 0)$  is an indicator function that obtains the value 1 if  $\|\mathbf{c}[i]\|_2 > 0$  and 0 otherwise, a block  $k$ -sparse vector  $\mathbf{c}$  can be defined by  $\|\mathbf{c}\|_{0,\mathcal{I}} \leq k$ .

Evidently, there is a one-to-one correspondence between a vector  $x$  in the union, and a block-sparse vector  $\mathbf{c}$ . The measurements (5) can also be represented explicitly in terms of  $\mathbf{c}$  as

$$\mathbf{y} = S^* x = S^* A\mathbf{c} = D\mathbf{c} \quad (16)$$

where  $\mathbf{D}$  is the  $n \times N$  matrix defined by

$$\mathbf{D} = \mathbf{S}^* \mathbf{A}. \quad (17)$$

We can therefore phrase our problem in terms of  $\mathbf{D}$  and  $\mathbf{c}$  as that of recovering a block- $k$  sparse vector  $\mathbf{c}$  over  $\mathcal{I}$  from the measurements (16).

Note that the choice of basis  $A_i$  for each subspace does not affect the sparsity pattern of our model. Indeed, choosing alternative bases will lead to  $x = \mathbf{A}\mathbf{W}\mathbf{c}$  where  $\mathbf{W}$  is a block diagonal matrix with blocks of size  $d_i$ . Defining  $\tilde{\mathbf{c}} = \mathbf{W}\mathbf{c}$ , the block sparsity pattern of  $\tilde{\mathbf{c}}$  is equal to that of  $\mathbf{c}$ . On the other hand, the basis can affect the properties of the resulting matrix  $\mathbf{D}$ . This point is beyond the scope of the present paper; further discussion can be found in [42].

Since our problem is equivalent to that of recovering a block sparse vector over  $\mathcal{I}$  from linear measurements  $\mathbf{y} = \mathbf{D}\mathbf{c}$ , in the remainder of the paper we focus our attention on this problem.

#### IV. UNIQUENESS AND STABILITY

In this section, we study the uniqueness and stability of our sampling method. These properties are intimately related to the RIP, which we generalize here to the block-sparse setting.

The first question we address is that of uniqueness, namely, conditions under which a block-sparse vector  $\mathbf{c}$  is uniquely determined by the measurement vector  $\mathbf{y} = \mathbf{D}\mathbf{c}$ .

*Proposition 1:* There is a unique block- $k$  sparse vector  $\mathbf{c}$  consistent with the measurements  $\mathbf{y} = \mathbf{D}\mathbf{c}$  if and only if  $\mathbf{D}\mathbf{c} \neq \mathbf{0}$  for every  $\mathbf{c} \neq \mathbf{0}$  that is block  $2k$ -sparse.

*Proof:* The proof follows from [24, Proposition 4].  $\square$

We next address the issue of stability. A sampling operator is stable for a set  $\mathcal{T}$  if and only if there exists constants  $\alpha > 0$ ,  $\beta < \infty$  such that

$$\alpha \|x_1 - x_2\|_{\mathcal{H}}^2 \leq \|S^*x_1 - S^*x_2\|_{\mathcal{H}}^2 \leq \beta \|x_1 - x_2\|_{\mathcal{H}}^2 \quad (18)$$

for every  $x_1, x_2$  in  $\mathcal{T}$ . The ratio  $\kappa = \beta/\alpha$  provides a measure for stability of the sampling operator. The operator is maximally stable when  $\kappa = 1$ . In our setting,  $S^*$  is replaced by  $\mathbf{D}$ , and the set  $\mathcal{T}$  contains block- $k$  sparse vectors. The next proposition follows immediately from (18) by noting that given two block- $k$  sparse vectors  $\mathbf{c}_1, \mathbf{c}_2$  their difference  $\mathbf{c}_1 - \mathbf{c}_2$  is block- $2k$  sparse.

*Proposition 2:* The measurement matrix  $\mathbf{D}$  is stable for every block  $k$ -sparse vector  $\mathbf{c}$  if and only if there exists  $C_1 > 0$  and  $C_2 < \infty$  such that

$$C_1 \|\mathbf{v}\|_2^2 \leq \|\mathbf{D}\mathbf{v}\|_2^2 \leq C_2 \|\mathbf{v}\|_2^2 \quad (19)$$

for every  $\mathbf{v}$  that is block  $2k$ -sparse.

It is easy to see that if  $\mathbf{D}$  satisfies (19) then  $\mathbf{D}\mathbf{c} \neq \mathbf{0}$  for all block  $2k$ -sparse vectors  $\mathbf{c}$ . Therefore, this condition implies both invertibility and stability.

##### A. Block RIP

Property (19) is related to the RIP used in several previous works in CS [9], [14], [15]. A matrix  $\mathbf{D}$  of size  $n \times N$  is said to have the RIP if there exists a constant  $\delta_k \in [0, 1)$  such that for every  $k$ -sparse  $\mathbf{c} \in \mathbb{R}^N$

$$(1 - \delta_k) \|\mathbf{c}\|_2^2 \leq \|\mathbf{D}\mathbf{c}\|_2^2 \leq (1 + \delta_k) \|\mathbf{c}\|_2^2. \quad (20)$$

Extending this property to block-sparse vectors leads to the following definition.

*Definition 2:* Let  $\mathbf{D} : \mathbb{R}^N \rightarrow \mathbb{R}^n$  be a given matrix. Then  $\mathbf{D}$  has the block RIP over  $\mathcal{I} = \{d_1, \dots, d_m\}$  with parameter  $\delta_{k|\mathcal{I}}$  if for every  $\mathbf{c} \in \mathbb{R}^N$  that is block  $k$ -sparse over  $\mathcal{I}$  we have that

$$(1 - \delta_{k|\mathcal{I}}) \|\mathbf{c}\|_2^2 \leq \|\mathbf{D}\mathbf{c}\|_2^2 \leq (1 + \delta_{k|\mathcal{I}}) \|\mathbf{c}\|_2^2. \quad (21)$$

By abuse of notation, we use  $\delta_k$  for the block-RIP constant  $\delta_{k|\mathcal{I}}$  when it is clear from the context that we refer to blocks. Block-RIP is a special case of the  $\mathcal{A}$ -restricted isometry defined in [25]. From Proposition 1, it follows that if  $\mathbf{D}$  satisfies the RIP (21) with  $\delta_{2k} < 1$ , then there is a unique block-sparse vector  $\mathbf{c}$  consistent with (16).

Note that a block  $k$ -sparse vector over  $\mathcal{I}$  is  $M$ -sparse in the conventional sense where  $M$  is the sum of the  $k$  largest values in  $\mathcal{I}$ , since it has at most  $M$  nonzero elements. If we require  $\mathbf{D}$  to satisfy RIP for all  $M$ -sparse vectors, then (21) must hold for all  $2M$ -sparse vectors  $\mathbf{c}$ . Since we only require the RIP for block sparse signals, (21) only has to be satisfied for a certain subset of  $2M$ -sparse signals, namely, those that have block sparsity. As a result, the block-RIP constant  $\delta_{k|\mathcal{I}}$  is typically smaller than  $\delta_M$  (where  $M$  depends on  $k$ ; for blocks with equal size  $d$ ,  $M = kd$ ).

To emphasize the advantage of block RIP over standard RIP, consider the following matrix, separated into three blocks of two columns each:

$$\mathbf{D} = \left( \begin{array}{cc|cc|cc} -1 & 1 & 0 & 0 & 0 & 1 \\ 0 & 2 & -1 & 0 & 0 & 3 \\ 0 & 3 & 0 & -1 & 0 & 1 \\ 0 & 1 & 0 & 0 & -1 & 1 \end{array} \right) \cdot \mathbf{B} \quad (22)$$

where  $\mathbf{B}$  is a diagonal matrix that results in  $\mathbf{D}$  with unit-norm columns, i.e.,  $\mathbf{B} = \text{diag}(1, 15, 1, 1, 1, 12)^{-1/2}$ . In this example,  $m = 3$  and  $\mathcal{I} = \{d_1 = 2, d_2 = 2, d_3 = 2\}$ . Suppose that  $\mathbf{c}$  is block-1 sparse, which corresponds to at most two nonzero values. Brute-force calculations show that the smallest value of  $\delta_2$  satisfying the standard RIP (20) is  $\delta_2 = 0.866$ . On the other hand, the block-RIP (21) corresponding to the case in which the two nonzero elements are restricted to occur in one block is satisfied with  $\delta_{1|\mathcal{I}} = 0.289$ . Increasing the number of nonzero elements to  $k = 4$ , we can verify that the standard RIP (20) does not hold for any  $\delta_4 \in [0, 1)$ . Indeed, in this example there exist two 4-sparse vectors that result in the same measurements. In contrast,  $\delta_{2|\mathcal{I}} = 0.966$  satisfies the lower bound in (21) when restricting the four nonzero values to two blocks. Consequently, the measurements  $\mathbf{y} = \mathbf{D}\mathbf{c}$  uniquely specify a single block-sparse  $\mathbf{c}$ .

In (5), we will see that the ability to recover  $\mathbf{c}$  in a computationally efficient way depends on the constant  $\delta_{2k|\mathcal{I}}$  in the block RIP (21). The smaller the value of  $\delta_{2k|\mathcal{I}}$ , the fewer samples are needed in order to guarantee stable recovery. Both standard and block RIP constants  $\delta_k, \delta_{k|\mathcal{I}}$  are by definition increasing with  $k$ . Therefore, it was suggested in [13] to normalize each of the columns of  $\mathbf{D}$  to 1, so as to start with  $\delta_1 = 0$ . In the same spirit, we recommend choosing the bases for  $\mathcal{A}_i$  such that  $\mathbf{D} = \mathbf{S}^* \mathbf{A}$  has unit-norm columns, corresponding to  $\delta_{1|\mathcal{I}} = 0$ .

*B. Recovery Method*

We have seen that if  $\mathbf{D}$  satisfies the RIP (21) with  $\delta_{2k} < 1$ , then there is a unique block-sparse vector  $\mathbf{c}$  consistent with (16). The question is how to find  $\mathbf{c}$  in practice. Below we present an algorithm that will in principle find the unique  $\mathbf{c}$  from the samples  $\mathbf{y}$ . Unfortunately, though, it has exponential complexity. In Section V, we show that under a stronger condition on  $\delta_{2k}$  we can recover  $\mathbf{c}$  in a stable and efficient manner.

Our first claim is that  $\mathbf{c}$  can be uniquely recovered by solving the optimization problem

$$\begin{aligned} \min_{\mathbf{c}} \quad & \|\mathbf{c}\|_{0,\mathcal{I}} \\ \text{s.t.} \quad & \mathbf{y} = \mathbf{D}\mathbf{c}. \end{aligned} \tag{23}$$

To show that (23) will indeed recover the true value of  $\mathbf{c}$ , suppose that there exists a  $\mathbf{c}'$  such that  $\mathbf{D}\mathbf{c}' = \mathbf{y}$  and  $\|\mathbf{c}'\|_{0,\mathcal{I}} \leq \|\mathbf{c}\|_{0,\mathcal{I}} \leq k$ . Since both  $\mathbf{c}, \mathbf{c}'$  are consistent with the measurements

$$\mathbf{0} = \mathbf{D}(\mathbf{c} - \mathbf{c}') = \mathbf{D}\mathbf{d} \tag{24}$$

where  $\|\mathbf{d}\|_{0,\mathcal{I}} \leq 2k$  so that  $\mathbf{d}$  is a block  $2k$ -sparse vector. Since  $\mathbf{D}$  satisfies (21) with  $\delta_{2k} < 1$ , we must have that  $\mathbf{d} = \mathbf{0}$  or  $\mathbf{c} = \mathbf{c}'$ .

In principle, (23) can be solved by searching over all possible sets of  $k$  blocks whether there exists a  $\mathbf{c}$  that is consistent with the measurements. The invertibility condition (21) ensures that there is only one such  $\mathbf{c}$ . However, clearly this approach is not efficient.

V. CONVEX RECOVERY ALGORITHM

A. Noise-Free Recovery

We now develop an efficient convex optimization problem instead of (23) to approximate  $\mathbf{c}$ . As we show, if  $\mathbf{D}$  satisfies (21) with a small enough value of  $\delta_{2k}$ , then the method we propose will recover  $\mathbf{c}$  exactly.

Our approach is to minimize the sum of the energy of the blocks  $\mathbf{c}[i]$ . To write down the problem explicitly, we define the mixed  $\ell_2/\ell_1$  norm over the index set  $\mathcal{I} = \{d_1, \dots, d_m\}$  as

$$\|\mathbf{c}\|_{2,\mathcal{I}} = \sum_{i=1}^m \|\mathbf{c}[i]\|_2. \tag{25}$$

The algorithm we suggest is then

$$\begin{aligned} \min_{\mathbf{c}} \quad & \|\mathbf{c}\|_{2,\mathcal{I}} \\ \text{s.t.} \quad & \mathbf{y} = \mathbf{D}\mathbf{c}. \end{aligned} \tag{26}$$

Problem (26) can be written as an SOCP by defining  $t_i = \|\mathbf{c}[i]\|_2$ . Then (26) is equivalent to

$$\begin{aligned} \min_{\mathbf{c}, t_i} \quad & \sum_{i=1}^m t_i \\ \text{s.t.} \quad & \mathbf{y} = \mathbf{D}\mathbf{c} \\ & t_i \geq \|\mathbf{c}[i]\|_2, \quad 1 \leq i \leq m \\ & t_i \geq 0, \quad 1 \leq i \leq m \end{aligned} \tag{27}$$

which can be solved using standard software packages.

The next theorem establishes that the solution to (26) is the true  $\mathbf{c}$  as long as  $\delta_{2k}$  is small enough.

*Theorem 1:* Let  $\mathbf{y} = \mathbf{D}\mathbf{c}_0$  be measurements of a block  $k$ -sparse vector  $\mathbf{c}_0$ . If  $\mathbf{D}$  satisfies the block RIP (21) with  $\delta_{2k} < \sqrt{2} - 1$  then

- 1) there is a unique block- $k$  sparse vector  $\mathbf{c}$  consistent with  $\mathbf{y}$ ;
- 2) the SOCP (27) has a unique solution;
- 3) the solution to the SOCP is equal to  $\mathbf{c}_0$ .

Before proving the theorem, we note that it provides a gain over standard CS results. Specifically, it is shown in [15] that if  $\mathbf{c}$  is  $k$ -sparse and the measurement matrix  $\mathbf{D}$  satisfies the standard RIP with  $\delta_{2k} < \sqrt{2} - 1$ , then  $\mathbf{c}$  can be recovered exactly from the measurements  $\mathbf{y} = \mathbf{D}\mathbf{c}$  via the linear program

$$\begin{aligned} \min_{\mathbf{c}} \quad & \|\mathbf{c}\|_1 \\ \text{s.t.} \quad & \mathbf{y} = \mathbf{D}\mathbf{c}. \end{aligned} \tag{28}$$

Since any block  $k$ -sparse vector is also  $M$ -sparse with  $M$  equal to the sum of the  $k$  largest values of  $d_i$ , we can find  $\mathbf{c}_0$  of Theorem 1 by solving (28) if  $\delta_{2M}$  is small enough. However, this standard CS approach does not exploit the fact that the nonzero values appear in blocks, and not in arbitrary locations within the vector  $\mathbf{c}_0$ . On the other hand, the SOCP (27) explicitly takes the block structure of  $\mathbf{c}_0$  into account. Therefore, the condition of Theorem 1 is not as stringent as that obtained by using equivalence results with respect to (28). Indeed, the block RIP (21) bounds the norm of  $\|\mathbf{D}\mathbf{c}\|$  over block sparse vectors  $\mathbf{c}$ , while the standard RIP considers all possible choices of  $\mathbf{c}$ , also those that are not  $2k$ -block sparse. Therefore, the value of  $\delta_{2k}$  in (21) can be lower than that obtained from (20) with  $k = 2M$ , as we illustrated by an example in Section III. This advantage will also be demonstrated by a concrete example at the end of the section.

Our proof below is rooted in that of [15]. However, some essential modifications are necessary in order to adapt the results to the block-sparse case. These differences are a result of the fact that our algorithm relies on the mixed  $\ell_2/\ell_1$  norm rather than the  $\ell_1$  norm alone. This adds another layer of complication to the proof, and therefore we expand the derivations in more detail than in [15].

*Proof of Theorem 1:* We first note that  $\delta_{2k} < 1$  guarantees uniqueness of  $\mathbf{c}_0$  from Proposition 1. To prove parts 2) and 3) we show that any solution to (26) has to be equal to  $\mathbf{c}_0$ . To this end, let  $\mathbf{c}' = \mathbf{c}_0 + \mathbf{h}$  be a solution of (26). The true value  $\mathbf{c}_0$  is nonzero over at most  $k$  blocks. We denote by  $\mathcal{I}_0$  the block indices for which  $\mathbf{c}_0$  is nonzero, and by  $\mathbf{h}_{\mathcal{I}_0}$  the restriction of  $\mathbf{h}$  to these blocks. Next, we decompose  $\mathbf{h}$  as

$$\mathbf{h} = \sum_{i=0}^{\ell-1} \mathbf{h}_{\mathcal{I}_i} \tag{29}$$

where  $\mathbf{h}_{\mathcal{I}_i}$  is the restriction of  $\mathbf{h}$  to the set  $\mathcal{I}_i$  which consists of  $k$  blocks, chosen such that the norm of  $\mathbf{h}_{\mathcal{I}_0}$  over  $\mathcal{I}_1$  is largest, the norm over  $\mathcal{I}_2$  is second largest, and so on. Our goal is to show that  $\mathbf{h} = \mathbf{0}$ . We prove this by noting that

$$\begin{aligned} \|\mathbf{h}\|_2 &= \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1} + \mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2 \\ &\leq \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 + \|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2. \end{aligned} \tag{30}$$

In the first part of the proof, we show that  $\|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2 \leq \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2$ . In the second part, we establish that  $\|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 = 0$ , which completes the proof.

*Part I:*  $\|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2 \leq \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2$ .

We begin by noting that

$$\|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2 = \left\| \sum_{i=2}^{\ell-1} \mathbf{h}_{\mathcal{I}_i} \right\|_2 \leq \sum_{i=2}^{\ell-1} \|\mathbf{h}_{\mathcal{I}_i}\|_2. \quad (31)$$

Therefore, it is sufficient to bound  $\|\mathbf{h}_{\mathcal{I}_i}\|_2$  for  $i \geq 2$ . Now

$$\|\mathbf{h}_{\mathcal{I}_i}\|_2 \leq k^{1/2} \|\mathbf{h}_{\mathcal{I}_i}\|_{\infty, \mathcal{I}} \leq k^{-1/2} \|\mathbf{h}_{\mathcal{I}_{i-1}}\|_{2, \mathcal{I}} \quad (32)$$

where we defined  $\|\mathbf{a}\|_{\infty, \mathcal{I}} = \max_i \|\mathbf{a}[i]\|_2$ . The first inequality follows from the fact that for any block  $k$ -sparse  $\mathbf{c}$

$$\|\mathbf{c}\|_2^2 = \sum_{|i|=k} \|\mathbf{c}[i]\|_2^2 \leq k \|\mathbf{c}\|_{\infty, \mathcal{I}}^2. \quad (33)$$

The second inequality in (32) is a result of the fact that the norm of each block in  $\mathbf{h}_{\mathcal{I}_i}$  is by definition smaller or equal to the norm of each block in  $\mathbf{h}_{\mathcal{I}_{i-1}}$ . Since there are at most  $k$  nonzero blocks,  $k \|\mathbf{h}_{\mathcal{I}_i}\|_{\infty, \mathcal{I}} \leq \|\mathbf{h}_{\mathcal{I}_{i-1}}\|_{2, \mathcal{I}}$ . Substituting (32) into (31)

$$\begin{aligned} \|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2 &\leq k^{-1/2} \sum_{i=1}^{\ell-2} \|\mathbf{h}_{\mathcal{I}_i}\|_{2, \mathcal{I}} \\ &\leq k^{-1/2} \sum_{i=1}^{\ell-1} \|\mathbf{h}_{\mathcal{I}_i}\|_{2, \mathcal{I}} \\ &= k^{-1/2} \|\mathbf{h}_{\mathcal{I}_0^c}\|_{2, \mathcal{I}} \end{aligned} \quad (34)$$

where the equality is a result of the fact that  $\|\mathbf{c}_1 + \mathbf{c}_2\|_{2, \mathcal{I}} = \|\mathbf{c}_1\|_{2, \mathcal{I}} + \|\mathbf{c}_2\|_{2, \mathcal{I}}$  if  $\mathbf{c}_1$  and  $\mathbf{c}_2$  are nonzero on disjoint blocks.

To develop a bound on  $\|\mathbf{h}_{\mathcal{I}_0^c}\|_{2, \mathcal{I}}$  note that since  $\mathbf{c}'$  is a solution to (26),  $\|\mathbf{c}_0\|_{2, \mathcal{I}} \geq \|\mathbf{c}'\|_{2, \mathcal{I}}$ . Using the fact that  $\mathbf{c}' = \mathbf{c}_0 + \mathbf{h}_{\mathcal{I}_0} + \mathbf{h}_{\mathcal{I}_0^c}$  and  $\mathbf{c}_0$  is supported on  $\mathcal{I}_0$  we have

$$\begin{aligned} \|\mathbf{c}_0\|_{2, \mathcal{I}} &\geq \|\mathbf{c}_0 + \mathbf{h}_{\mathcal{I}_0}\|_{2, \mathcal{I}} + \|\mathbf{h}_{\mathcal{I}_0^c}\|_{2, \mathcal{I}} \\ &\geq \|\mathbf{c}_0\|_{2, \mathcal{I}} - \|\mathbf{h}_{\mathcal{I}_0}\|_{2, \mathcal{I}} + \|\mathbf{h}_{\mathcal{I}_0^c}\|_{2, \mathcal{I}} \end{aligned} \quad (35)$$

from which we conclude that

$$\|\mathbf{h}_{\mathcal{I}_0^c}\|_{2, \mathcal{I}} \leq \|\mathbf{h}_{\mathcal{I}_0}\|_{2, \mathcal{I}} \leq k^{1/2} \|\mathbf{h}_{\mathcal{I}_0}\|_2. \quad (36)$$

The last inequality follows from applying Cauchy–Schwarz to any block  $k$ -sparse vector  $\mathbf{c}$

$$\|\mathbf{c}\|_{2, \mathcal{I}} = \sum_{|i|=k} \|\mathbf{c}[i]\|_2 \cdot 1 \leq k^{1/2} \|\mathbf{c}\|_2. \quad (37)$$

Substituting (36) into (34)

$$\|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2 \leq \|\mathbf{h}_{\mathcal{I}_0}\|_2 \leq \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 \quad (38)$$

which completes the first part of the proof.

*Part II:*  $\|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 = 0$ .

We next show that  $\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}$  must be equal to  $\mathbf{0}$ . In this part we invoke the RIP.

Since  $D\mathbf{c}_0 = D\mathbf{c}' = \mathbf{y}$ , we have  $D\mathbf{h} = \mathbf{0}$ . Using the fact that  $\mathbf{h} = \mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1} + \sum_{i \geq 2} \mathbf{h}_{\mathcal{I}_i}$

$$\|D\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 = - \sum_{i=2}^{\ell-1} \langle D(\mathbf{h}_{\mathcal{I}_0} + \mathbf{h}_{\mathcal{I}_1}), D\mathbf{h}_{\mathcal{I}_i} \rangle. \quad (39)$$

From the parallelogram identity and the block-RIP it can be shown that

$$|\langle D\mathbf{c}_1, D\mathbf{c}_2 \rangle| \leq \delta_{2k} \|\mathbf{c}_1\|_2 \|\mathbf{c}_2\|_2 \quad (40)$$

for any two-block  $k$ -sparse vectors with disjoint support. The proof is similar to [15, Lemma 2.1] for the standard RIP. Therefore

$$|\langle D\mathbf{h}_{\mathcal{I}_0}, D\mathbf{h}_{\mathcal{I}_i} \rangle| \leq \delta_{2k} \|\mathbf{h}_{\mathcal{I}_0}\|_2 \|\mathbf{h}_{\mathcal{I}_i}\|_2 \quad (41)$$

and similarly for  $\langle D\mathbf{h}_{\mathcal{I}_1}, D\mathbf{h}_{\mathcal{I}_i} \rangle$ . Substituting into (39)

$$\begin{aligned} \|D\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 &= \left| \sum_{i=2}^{\ell-1} \langle D(\mathbf{h}_{\mathcal{I}_0} + \mathbf{h}_{\mathcal{I}_1}), D\mathbf{h}_{\mathcal{I}_i} \rangle \right| \\ &\leq \sum_{i=2}^{\ell-1} (|\langle D\mathbf{h}_{\mathcal{I}_0}, D\mathbf{h}_{\mathcal{I}_i} \rangle| + |\langle D\mathbf{h}_{\mathcal{I}_1}, D\mathbf{h}_{\mathcal{I}_i} \rangle|) \\ &\leq \delta_{2k} (\|\mathbf{h}_{\mathcal{I}_0}\|_2 + \|\mathbf{h}_{\mathcal{I}_1}\|_2) \sum_{i=2}^{\ell-1} \|\mathbf{h}_{\mathcal{I}_i}\|_2. \end{aligned} \quad (42)$$

From the Cauchy–Schwarz inequality, any length-2 vector  $\mathbf{a}$  satisfies  $\mathbf{a}(1) + \mathbf{a}(2) \leq \sqrt{2} \|\mathbf{a}\|_2$ . Therefore

$$\begin{aligned} \|\mathbf{h}_{\mathcal{I}_0}\|_2 + \|\mathbf{h}_{\mathcal{I}_1}\|_2 &\leq \sqrt{2} \sqrt{\|\mathbf{h}_{\mathcal{I}_0}\|_2^2 + \|\mathbf{h}_{\mathcal{I}_1}\|_2^2} \\ &= \sqrt{2} \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 \end{aligned} \quad (43)$$

where the last equality is a result of the fact that  $\mathbf{h}_{\mathcal{I}_0}$  and  $\mathbf{h}_{\mathcal{I}_1}$  have disjoint support. Substituting into (42) and using (32), (34), and (36)

$$\begin{aligned} \|D\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 &\stackrel{(32), (34)}{\leq} \sqrt{2} k^{-1/2} \delta_{2k} \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 \|\mathbf{h}_{\mathcal{I}_0^c}\|_{2, \mathcal{I}} \\ &\stackrel{(36)}{\leq} \sqrt{2} \delta_{2k} \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 \|\mathbf{h}_{\mathcal{I}_0}\|_2 \\ &\leq \sqrt{2} \delta_{2k} \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 \end{aligned} \quad (44)$$

where the last inequality follows from  $\|\mathbf{h}_{\mathcal{I}_0}\|_2 \leq \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2$ . Combining (44) with the RIP (21) we have

$$(1 - \delta_{2k}) \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 \leq \|D\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 \leq \sqrt{2} \delta_{2k} \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2. \quad (45)$$

Since  $\delta_{2k} < \sqrt{2} - 1$ , (45) can hold only if  $\|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 = 0$ , which completes the proof.  $\square$

We conclude this subsection by pointing out more explicitly the differences between the proof of Theorem 1 and that in [15]. The main difference begins in (32); in our formulation, each of the subvectors  $\mathbf{h}_{\mathcal{I}_i}$  may have a different number of nonzero elements, while the equivalent equation in [15] (10) relies on the fact that the maximal number of nonzero elements in each of the subvectors is the same. This requires the use of several mixed

norms in our setting. The rest of the proof follows the spirit of [15], where in some of the inequalities conventional norms are used, while in others the adaptation to our setting necessitates mixed norms.

### B. Robust Recovery

We now treat the situation in which the observations are noisy, and the vector  $\mathbf{c}_0$  is not exactly block- $k$  sparse.

Specifically, suppose that the measurements (16) are corrupted by bounded noise so that

$$\mathbf{y} = \mathbf{D}\mathbf{c} + \mathbf{z} \quad (46)$$

where  $\|\mathbf{z}\|_2 \leq \epsilon$ . In order to recover  $\mathbf{c}$ , we use the modified SOCP

$$\begin{aligned} \min_{\mathbf{c}} \quad & \|\mathbf{c}\|_{2,\mathcal{I}} \\ \text{s.t.} \quad & \|\mathbf{y} - \mathbf{D}\mathbf{c}\|_2 \leq \epsilon. \end{aligned} \quad (47)$$

In addition, given a  $\mathbf{c} \in \mathbb{R}^N$ , we denote by  $\mathbf{c}^k$  the best approximation of  $\mathbf{c}$  by a vector with  $k$  nonzero blocks, so that  $\mathbf{c}^k$  minimizes  $\|\mathbf{c} - \mathbf{d}\|_{2,\mathcal{I}}$  over all block  $k$ -sparse vectors  $\mathbf{d}$ . Theorem 2 shows that even when  $\mathbf{c}$  is not block  $k$ -sparse and the measurements are noisy, the best block- $k$  approximation can be well approximated using (47).

*Theorem 2:* Let  $\mathbf{y} = \mathbf{D}\mathbf{c}_0 + \mathbf{z}$  be noisy measurements of a vector  $\mathbf{c}_0$ . Let  $\mathbf{c}^k$  denote the best block  $k$ -sparse approximation of  $\mathbf{c}_0$ , such that  $\mathbf{c}^k$  is block  $k$ -sparse and minimizes  $\|\mathbf{c}_0 - \mathbf{d}\|_{2,\mathcal{I}}$  over all block  $k$ -sparse vectors  $\mathbf{d}$ , and let  $\mathbf{c}'$  be a solution to (47). If  $\mathbf{D}$  satisfies the block RIP (21) with  $\delta_{2k} < \sqrt{2} - 1$  then

$$\begin{aligned} \|\mathbf{c}_0 - \mathbf{c}'\|_2 \leq & \frac{2(1 - \delta_{2k})}{1 - (1 + \sqrt{2})\delta_{2k}} k^{-1/2} \|\mathbf{c}_0 - \mathbf{c}^k\|_{2,\mathcal{I}} \\ & + \frac{4\sqrt{1 + \delta_{2k}}}{1 - (1 + \sqrt{2})\delta_{2k}} \epsilon. \end{aligned} \quad (48)$$

Before proving the theorem, note that the first term in (48) is a result of the fact that  $\mathbf{c}_0$  is not exactly  $k$ -block sparse. The second expression quantifies the recovery error due to the noise.

*Proof:* The proof is very similar to that of Theorem 1 with a few differences which we indicate. These changes follow the proof of [15, Theorem 1.3], with appropriate modifications to address the mixed norm.

Denote by  $\mathbf{c}' = \mathbf{c}_0 + \mathbf{h}$  the solution to (47). Due to the noise and the fact that  $\mathbf{c}_0$  is not block  $k$ -sparse, we will no longer obtain  $\mathbf{h} = \mathbf{0}$ . However, we will show that  $\|\mathbf{h}\|_2$  is bounded. To this end, we begin as in the proof of Theorem 1 by using (30). In the first part of the proof we show that  $\|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2 \leq \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 + 2e_0$  where  $e_0 = k^{-1/2} \|\mathbf{c}_0 - \mathbf{c}_{\mathcal{I}_0}\|_{2,\mathcal{I}}$  and  $\mathbf{c}_{\mathcal{I}_0}$  is the restriction of  $\mathbf{c}_0$  onto the  $k$  blocks corresponding to the largest  $\ell_2$  norm. Note that  $\mathbf{c}_{\mathcal{I}_0} = \mathbf{c}^k$ . In the second part, we develop a bound on  $\|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2$ .

*Part I: Bound on  $\|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2$ .*

We begin by decomposing  $\mathbf{h}$  as in the proof of Theorem 1. The inequalities until (35) hold here as well. Instead of (35) we have

$$\begin{aligned} \|\mathbf{c}_0\|_{2,\mathcal{I}} & \geq \|\mathbf{c}_{\mathcal{I}_0} + \mathbf{h}_{\mathcal{I}_0}\|_{2,\mathcal{I}} + \|\mathbf{c}_{\mathcal{I}_0^c} + \mathbf{h}_{\mathcal{I}_0^c}\|_{2,\mathcal{I}} \\ & \geq \|\mathbf{c}_{\mathcal{I}_0}\|_{2,\mathcal{I}} - \|\mathbf{h}_{\mathcal{I}_0}\|_{2,\mathcal{I}} + \|\mathbf{h}_{\mathcal{I}_0^c}\|_{2,\mathcal{I}} - \|\mathbf{c}_{\mathcal{I}_0^c}\|_{2,\mathcal{I}}. \end{aligned} \quad (49)$$

Therefore

$$\|\mathbf{h}_{\mathcal{I}_0^c}\|_{2,\mathcal{I}} \leq 2\|\mathbf{c}_{\mathcal{I}_0^c}\|_{2,\mathcal{I}} + \|\mathbf{h}_{\mathcal{I}_0}\|_{2,\mathcal{I}} \quad (50)$$

where we used the fact that  $\|\mathbf{c}_0\|_{2,\mathcal{I}} - \|\mathbf{c}_{\mathcal{I}_0}\|_{2,\mathcal{I}} = \|\mathbf{c}_{\mathcal{I}_0^c}\|_{2,\mathcal{I}}$ . Combining (34), (37), and (50) we have

$$\|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2 \leq \|\mathbf{h}_{\mathcal{I}_0}\|_2 + 2e_0 \leq \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 + 2e_0 \quad (51)$$

where  $e_0 = k^{-1/2} \|\mathbf{c}_0 - \mathbf{c}_{\mathcal{I}_0}\|_{2,\mathcal{I}}$ .

*Part II: Bound on  $\|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2$ .*

Using the fact that  $\mathbf{h} = \mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1} + \sum_{i \geq 2} \mathbf{h}_{\mathcal{I}_i}$ , we have

$$\begin{aligned} \|\mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 & = \langle \mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}, \mathbf{D}\mathbf{h} \rangle - \sum_{i=2}^{\ell-1} \langle \mathbf{D}(\mathbf{h}_{\mathcal{I}_0} + \mathbf{h}_{\mathcal{I}_1}), \mathbf{D}\mathbf{h}_{\mathcal{I}_i} \rangle. \end{aligned} \quad (52)$$

From (21)

$$\begin{aligned} |\langle \mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}, \mathbf{D}\mathbf{h} \rangle| & \leq \|\mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 \|\mathbf{D}\mathbf{h}\|_2 \\ & \leq \sqrt{1 + \delta_{2k}} \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 \|\mathbf{D}\mathbf{h}\|_2. \end{aligned} \quad (53)$$

Since both  $\mathbf{c}'$  and  $\mathbf{c}_0$  are feasible

$$\begin{aligned} \|\mathbf{D}\mathbf{h}\|_2 = \|\mathbf{D}(\mathbf{c}_0 - \mathbf{c}')\|_2 & \leq \|\mathbf{D}\mathbf{c}_0 - \mathbf{y}\|_2 + \|\mathbf{D}\mathbf{c}' - \mathbf{y}\|_2 \\ & \leq 2\epsilon \end{aligned} \quad (54)$$

and (53) becomes

$$|\langle \mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}, \mathbf{D}\mathbf{h} \rangle| \leq 2\epsilon \sqrt{1 + \delta_{2k}} \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2. \quad (55)$$

Substituting into (52)

$$\begin{aligned} \|\mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 & \leq |\langle \mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}, \mathbf{D}\mathbf{h} \rangle| + \sum_{i=2}^{\ell-1} |\langle \mathbf{D}(\mathbf{h}_{\mathcal{I}_0} + \mathbf{h}_{\mathcal{I}_1}), \mathbf{D}\mathbf{h}_{\mathcal{I}_i} \rangle| \\ & \leq 2\epsilon \sqrt{1 + \delta_{2k}} \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 + \sum_{i=2}^{\ell-1} |\langle \mathbf{D}(\mathbf{h}_{\mathcal{I}_0} + \mathbf{h}_{\mathcal{I}_1}), \mathbf{D}\mathbf{h}_{\mathcal{I}_i} \rangle|. \end{aligned} \quad (56)$$

Combining with (42) and (44)

$$\begin{aligned} \|\mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 & \leq \left( 2\epsilon \sqrt{1 + \delta_{2k}} + \sqrt{2}\delta_{2k} k^{-1/2} \|\mathbf{h}_{\mathcal{I}_0^c}\|_{2,\mathcal{I}} \right) \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2. \end{aligned} \quad (57)$$

Using (37) and (50) we have the upper bound

$$\begin{aligned} \|\mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 & \leq \left( 2\epsilon \sqrt{1 + \delta_{2k}} + \sqrt{2}\delta_{2k} (\|\mathbf{h}_{\mathcal{I}_0}\| + 2e_0) \right) \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2. \end{aligned} \quad (58)$$

On the other hand, the RIP results in the lower bound

$$\|\mathbf{D}\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2 \geq (1 - \delta_{2k}) \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2^2. \quad (59)$$

From (58) and (59),

$$\begin{aligned} (1 - \delta_{2k}) \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 & \leq 2\epsilon \sqrt{1 + \delta_{2k}} + \sqrt{2}\delta_{2k} (\|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\| + 2e_0) \end{aligned} \quad (60)$$

or

$$\|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 \leq \frac{2\sqrt{1 + \delta_{2k}}}{1 - (1 + \sqrt{2})\delta_{2k}} \epsilon + \frac{2\sqrt{2}\delta_{2k}}{1 - (1 + \sqrt{2})\delta_{2k}} e_0. \quad (61)$$



TABLE I  
COMPARISON OF ALGORITHMS FOR SIGNAL RECOVERY FROM  $\mathbf{y} = \mathbf{D}\mathbf{c}_0 + \mathbf{z}$

	Algorithm (26)	Algorithm (47)
$\mathbf{c}_0$	block $k$ -sparse	arbitrary
Noise $\mathbf{z}$	none ( $\mathbf{z} = \mathbf{0}$ )	bounded $\ \mathbf{z}\ _2 \leq \epsilon$
Condition on $\mathbf{D}$	$\delta_{2k} \leq \sqrt{2} - 1$	$\delta_{2k} \leq \sqrt{2} - 1$
Recovery $\mathbf{c}'$	$\mathbf{c}' = \mathbf{c}_0$	$\ \mathbf{c}_0 - \mathbf{c}'\ _2$ small; see (48)

The condition  $\delta_{2k} < \sqrt{2} - 1$  ensures that the denominator in (61) is positive. Substituting (61) results in

$$\|\mathbf{h}\|_2 \leq \|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 + \|\mathbf{h}_{(\mathcal{I}_0 \cup \mathcal{I}_1)^c}\|_2 \leq 2\|\mathbf{h}_{\mathcal{I}_0 \cup \mathcal{I}_1}\|_2 + 2\epsilon_0 \quad (62)$$

which completes the proof of the theorem.  $\square$

To summarize this section, we have seen that as long as  $\mathbf{D}$  satisfies the block-RIP (21) with a suitable constant, any block- $k$  sparse vector can be perfectly recovered from its samples  $\mathbf{y} = \mathbf{D}\mathbf{c}$  using the convex SOCP (26). This algorithm is stable in the sense that by slightly modifying it as in (47) it can tolerate noise in a way that ensures that the norm of the recovery error is bounded by the noise level. Furthermore, if  $\mathbf{c}$  is not block  $k$ -sparse, then its best block-sparse approximation can be approached by solving the SOCP. These results are summarized in Table I. In the table,  $\delta_{2k}$  refers to the block RIP constant.

### C. Advantage of Block Sparsity

The standard sparsity model considered in CS assumes that  $\mathbf{x}$  has at most  $k$  nonzero elements, however it does not impose any further structure. In particular, the nonzero components can appear anywhere in the vector. There are many practical scenarios in which the nonzero values are aligned to blocks, meaning they appear in regions, and are not arbitrarily spread throughout the vector. One example in the structured union of subspaces model we treat in this paper. Other examples are considered in [27].

Prior work on recovery of block-sparse vectors [26] assumed consecutive blocks of the same size. It was shown that in this case, when  $n, N$  go to infinity, the algorithm (26) will recover the true block-sparse vector with overwhelming probability. Their analysis is based on characterization of the null space of  $\mathbf{D}$ . In contrast, our approach relies on RIP which allows the derivation of uniqueness and equivalence conditions for finite dimensions and not only in the asymptotic regime. In addition, Theorem 2 considers the case of mismodeling and noisy observations while in [26] only the ideal noise-free setting is treated.

To demonstrate the advantage of our algorithm over standard basis pursuit (28), consider the matrix  $\mathbf{D}$  of (22). In Section V, the standard and block RIP constants of  $\mathbf{D}$  were calculated and it was shown that block RIP constants are smaller. This suggests that there are input vectors  $\mathbf{x}$  for which the mixed  $\ell_2/\ell_1$  method of (26) will be able to recover them exactly from measurements  $\mathbf{y} = \mathbf{D}\mathbf{c}$  while standard  $\ell_1$  minimization will fail. To illustrate this behavior, let  $\mathbf{x} = [0, 0, 1, -1, -1, 0.1]^T$  be a 4-sparse vector, in which the nonzero elements are known to appear in blocks of length 2. The prior knowledge that  $\mathbf{x}$  is 4-sparse is not sufficient to determine  $\mathbf{x}$  from  $\mathbf{y}$ . In contrast, there is a unique block-sparse vector consistent with  $\mathbf{y}$ . Furthermore, our algorithm which is a relaxed version of (23), finds the correct  $\mathbf{x}$  while standard  $\ell_1$  minimization fails in this case; its output is  $\hat{\mathbf{x}} = [-0.0289, 0, 0.9134, -1.0289, -1.0289, 0]$ .

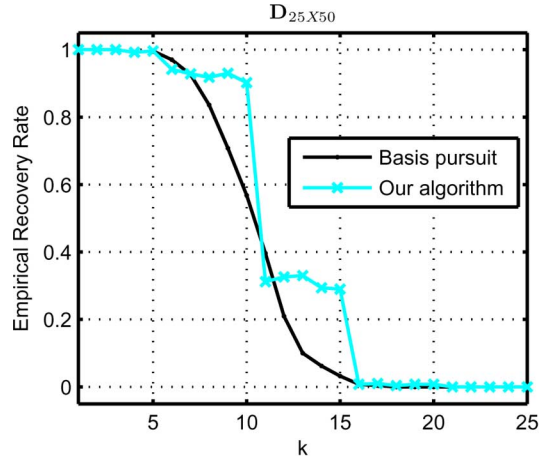


Fig. 2. Recovery rate of block-sparse signals using standard  $\ell_1$  minimization (basis pursuit) and the mixed  $\ell_2/\ell_1$  algorithm.

We further compare the recovery performance of  $\ell_1$  minimization (28) and our algorithm (26) for an extensive set of random signals. In the experiment, we draw a matrix  $\mathbf{D}$  of size  $25 \times 50$  from the Gaussian ensemble. The input vector  $\mathbf{x}$  is also randomly generated as a block-sparse vector with blocks of length 5. We draw  $1 \leq k \leq 25$  nonzero entries from a zero-mean unit variance normal distribution and divide them into blocks which are chosen uniformly at random within  $\mathbf{x}$ . Each of the algorithms is executed based on the measurements  $\mathbf{y} = \mathbf{D}\mathbf{x}$ . In Fig. 2, we plot the fraction of successful reconstructions for each  $k$  over 500 experiments. The results illustrate the advantage of incorporating the block-sparsity structure into the optimization program. An interesting feature of the graph is that when using the block-sparse recovery approach, the performance is roughly constant over the block length (5 in this example). This explains the performance advantage over standard sparse recovery.

## VI. APPLICATION TO MMV MODELS

We now specialize our algorithm and equivalence results to the MMV problem. This leads to two contributions which we discuss in this section: The first is an equivalence result based on RIP for a mixed-norm MMV algorithm. The second is a new measurement strategy in MMV problems that leads to improved performance over conventional MMV methods, both in simulations and as measured by the RIP-based equivalence condition. In contrast to previous equivalence results, for this strategy we show that even if we choose the worst possible  $\mathbf{X}$ , improved performance over the single measurement setting can be guaranteed.

### A. Equivalence Results

As we have seen in Section II, a special case of block sparsity is the MMV model, in which we are given a matrix of measurements  $\mathbf{Y} = \mathbf{M}\mathbf{X}$  where  $\mathbf{X}$  is an unknown  $L \times d$  matrix that has at most  $k$  nonzero rows. Denoting by  $\mathbf{c} = \text{vec}(\mathbf{X}^T)$ ,  $\mathbf{y} = \text{vec}(\mathbf{Y}^T)$ ,  $\mathbf{D} = \mathbf{M}^T \otimes \mathbf{I}_d$  we can express the vector of measurements  $\mathbf{y}$  as  $\mathbf{y} = \mathbf{D}\mathbf{c}$  where  $\mathbf{c}$  is a block sparse vector with consecutive blocks of length  $d$ . Therefore, the results of Theorems 1 and 2 can be specified to this problem.

Recovery algorithms for MMV using convex optimization programs were studied in [30], [32] and several greedy algorithms were proposed in [29], [31]. Specifically, in [29]–[32], the authors study a class of optimization programs, which we refer to as M-BP

$$\text{M-BP}(\ell_q) : \min \sum_{i=1}^L \|\mathbf{X}^i\|_q^p \quad \text{s.t. } \mathbf{Y} = \mathbf{M}\mathbf{X} \quad (63)$$

where  $\mathbf{X}^i$  is the  $i$ th row of  $\mathbf{X}$ . The choice  $p = 1, q = \infty$  was considered in [32], while [30] treated the case of  $p = 1$  and arbitrary  $q$ . Using  $p \leq 1$  and  $q = 2$  was suggested in [29], [43], leading to the iterative algorithm M-FOCUSS. For  $p = 1, q = 2$ , the program (63) has a global minimum which M-FOCUSS is proven to find. A nice comparison between these methods can be found in [32]. Equivalence for MMV algorithms based on RIP analysis does not appear in previous papers. The most detailed theoretical analysis can be found in [30] which establishes equivalence results based on mutual coherence. The results imply equivalence for (63) with  $p = 1$  under conditions equal to those obtained for the single measurement case. Note that RIP analysis typically leads to tighter equivalence bounds than mutual coherence analysis.

In our recent work [20], we suggested an alternative approach to solving MMV problems. Our strategy is based on merging the  $d$  measurement columns with random coefficients and in such a way transforming the multiple measurement problem into a single measurement counterpart. As proved in [20], this technique preserves the nonzero location set with probability one thus reducing computational complexity. Moreover, we showed that this method can be used to boost the empirical recovery rate by repeating the random merging several times.

Using the block-sparsity approach we can alternatively cast any MMV model as a single measurement vector problem by deterministically transforming the multiple measurement vectors into the single vector model  $\text{vec}(\mathbf{Y}^T) = (\mathbf{M} \otimes \mathbf{I}_d)\text{vec}(\mathbf{X}^T)$ . Here,  $\mathbf{c} = \text{vec}(\mathbf{X}^T)$  is block- $k$  sparse with consecutive blocks of length  $d$ . In contrast to [20], this does not reduce the number of unknowns so that the computational complexity of the resulting algorithm is on the same order as previous approaches, and also does not offer the opportunity for boosting. However, as seen in the next subsection, with an appropriate choice of measurement matrix this approach results in improved recovery capabilities.

Since we can cast the MMV problem as one of block-sparse recovery, we may apply our equivalence results of Theorem 1 to this setting leading to RIP-based equivalence. To this end, we first note that applying the SOCP (26) to the effective measurement vector  $\mathbf{y}$  is the same as solving (63) with  $p = 1, q = 2$ . Thus, the equivalence conditions we develop below relate to this program. Next, if  $\mathbf{z} = \mathbf{D}\mathbf{c}$  where  $\mathbf{c}$  is a block  $2k$ -sparse vector and  $\mathbf{D} = \mathbf{M} \otimes \mathbf{I}_d$ , then taking the structure of  $\mathbf{D}$  into account,  $\mathbf{Z} = \mathbf{M}\mathbf{X}$  where  $\mathbf{X}$  is a size  $L \times d$  matrix whose  $i$ th row is equal to  $\mathbf{c}[i]$ , and similarly for  $\mathbf{Z}$ . The block sparsity of  $\mathbf{c}$  implies that  $\mathbf{X}$  has at most  $2k$  nonzero rows. The squared  $\ell_2$  norm  $\|\mathbf{z}\|_2^2$  is equal to the squared  $\ell_2$  norm of the rows of  $\mathbf{Z}$  which can be written as

$$\|\mathbf{z}\|_2^2 = \|\mathbf{Z}\|_F^2 = \text{Tr}(\mathbf{Z}^T \mathbf{Z}) \quad (64)$$

where  $\|\mathbf{Z}\|_F$  denotes the Frobenius norm. Since  $\|\mathbf{c}\|_2^2 = \|\mathbf{X}\|_F^2$  the RIP condition becomes

$$(1 - \delta_{2k})\text{Tr}(\mathbf{X}^T \mathbf{X}) \leq \text{Tr}(\mathbf{X}^T \mathbf{M}^T \mathbf{M} \mathbf{X}) \leq (1 + \delta_{2k})\text{Tr}(\mathbf{X}^T \mathbf{X}) \quad (65)$$

for any  $L \times d$  matrix  $\mathbf{X}$  with at most  $2k$  nonzero rows.

We now show that (65) is equivalent to the standard RIP condition

$$(1 - \delta_{2k})\|\mathbf{x}\|_2^2 \leq \|\mathbf{M}\mathbf{x}\|_2^2 \leq (1 + \delta_{2k})\|\mathbf{x}\|_2^2 \quad (66)$$

for any length  $L$  vector  $\mathbf{x}$  that is  $2k$ -sparse. To see this, suppose first that (65) is satisfied for every matrix  $\mathbf{X}$  with at most  $2k$  nonzero rows and let  $\mathbf{x}$  be an arbitrary  $2k$ -sparse vector. If we define  $\mathbf{X}$  to be the matrix whose columns are all equal to  $\mathbf{x}$ , then  $\mathbf{X}$  will have at most  $2k$  nonzero rows and therefore satisfies (65). Since the columns of  $\mathbf{X}$  are all equal,  $\text{Tr}(\mathbf{X}^T \mathbf{X}) = d\|\mathbf{x}\|_2^2$  and  $\text{Tr}(\mathbf{X}^T \mathbf{M}^T \mathbf{M} \mathbf{X}) = d\|\mathbf{M}\mathbf{x}\|_2^2$  so that (66) holds. Conversely, suppose that (66) is satisfied for all  $2k$ -sparse vectors  $\mathbf{x}$  and let  $\mathbf{X}$  be an arbitrary matrix with at most  $2k$  nonzero rows. Denoting by  $\mathbf{x}_j$  the columns of  $\mathbf{X}$ , each  $\mathbf{x}_j$  is  $2k$ -sparse and therefore satisfies (66). Summing over all values  $j$  results in (65).

To summarize, if  $\mathbf{M}$  satisfies the conventional RIP condition (66), then the algorithm (63) with  $p = 1, q = 2$  will recover the true unknown  $\mathbf{X}$ . This requirement reduces to that we would obtain if we tried to recover each column of  $\mathbf{X}$  separately, using the standard  $\ell_1$  approach (28). As we already noted, previous equivalence results for MMV algorithms also share this feature. Although this condition guarantees that processing the vectors jointly does not harm the recovery ability, in practice exploiting the joint sparsity pattern of  $\mathbf{X}$  via (63) leads to improved results. Unfortunately, this behavior is not captured by any of the known equivalence conditions. This is due to the special structure of  $\mathbf{D} = \mathbf{M} \otimes \mathbf{I}$ . Since each measurement vector  $\mathbf{y}_i$  is affected only by the corresponding vector  $\mathbf{x}_i$ , it is clear that in the worst case we can choose  $\mathbf{x}_i = \mathbf{x}$  for some vector  $\mathbf{x}$ . In this case, all the  $\mathbf{y}_i$ 's are equal so that adding measurement vectors will not improve our recovery ability. Consequently, worst case analysis based on the standard measurement model for MMV problems cannot lead to improved performance over the single measurement case.

### B. Improved MMV Recovery

We have seen that the pessimistic equivalence results for MMV algorithms is a consequence of the fact that in the worst case scenario in which  $\mathbf{x}_i = \mathbf{x}$ , using a separable measurement strategy will render all observation vectors equal. In this subsection, we introduce an alternative measurement technique for MMV problems that can lead to improved worst case behavior, as measured by RIP, over the single channel case.

One way to improve the analytical results is to consider an average case analysis instead of a worst case approach. In [44], we show that if the unknown vectors  $\mathbf{x}_i$  are generated randomly, then the performance improves with increasing number of measurement vectors. The advantage stems from the fact that the situation of equal vectors has zero probability and therefore does not affect the average performance. Here we take a different route which does not involve randomness in the unknown

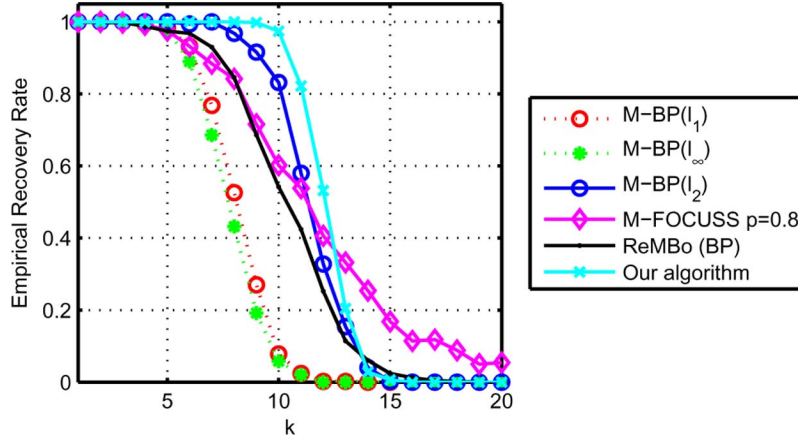


Fig. 3. Recovery rate for different number  $k$  of nonzero rows in  $\mathbf{X}$ . Each point on the graph represents an average recovery rate over 500 simulations.

vectors, and leads to improved results even in the worst case (namely, without requiring an average analysis).

To enhance the performance of MMV recovery, we note that when we allow for an arbitrary (unstructured)  $\mathbf{D}$ , the RIP condition of Theorem 1 is weaker than the standard RIP requirement for recovering  $k$ -sparse vectors. This suggests that we can improve the performance of MMV methods by converting the problem into a general block sparsity problem, and then sampling with an arbitrary unstructured matrix  $\mathbf{D}$  rather than the choice  $\mathbf{D} = \mathbf{M}^T \otimes \mathbf{I}_d$ . The tradeoff introduced is increased computational complexity since each measurement is based on all input vectors. The theoretical conditions will now be looser, since block-RIP is weaker than standard RIP. Furthermore, in practice, this approach often improves the performance over separable MMV measurement techniques as we illustrate in the following example.

In the example, we compare the performance of several MMV algorithms for recovering  $\mathbf{X}$  in the model  $\mathbf{Y} = \mathbf{M}\mathbf{X}$ , with our method based on block sparsity in which the measurements  $\mathbf{y}$  are obtained via  $\mathbf{y} = \mathbf{D}\mathbf{c}$  where  $\mathbf{c} = \text{vec}(\mathbf{X}^T)$  and  $\mathbf{D}$  is a dense matrix. Choosing  $\mathbf{D}$  as a block diagonal matrix with blocks equal to  $\mathbf{M}$  results in the standard MMV measurement model. The effective matrices  $\mathbf{D}$  have the same size in the case in which it is block diagonal and when it is dense. To compare the performance of (26) with a dense  $\mathbf{D}$  to that of (63) with a block diagonal  $\mathbf{D}$ , we compute the empirical recovery rate of the methods in the same way performed in [20]. The matrices  $\mathbf{M}$  and  $\mathbf{D}$  are drawn randomly from a Gaussian ensemble. In our example, we choose  $\ell = 20$ ,  $L = 30$ ,  $d = 5$  where  $\ell$  is the number of rows in  $\mathbf{Y}$ . The matrix  $\mathbf{X}$  is generated randomly by first selecting the  $k$  nonzero rows uniformly at random, and then drawing the elements in these rows from a normal distribution. The empirical recovery rates using the methods of (63) for different choices of  $q$  and  $p$ , ReMBO [20] and our algorithm (26) with dense  $\mathbf{D}$  are depicted in Fig. 3. When the index  $p$  is omitted it is equal to 1. Evidently, our algorithm performs better than most popular optimization techniques for MMV systems. We stress that the performance advantage is due to the joint measurement process rather than a new recovery algorithm.

## VII. RANDOM MATRICES

Theorems 1 and 2 establish that a sufficiently small block RIP constant  $\delta_{2k|\mathcal{I}}$  ensures exact recovery of the coefficient vector  $\mathbf{c}$ . We now prove that random matrices are likely to satisfy this requirement. Specifically, we show that the probability that  $\delta_{k|\mathcal{I}}$  exceeds a certain threshold decays exponentially in the length of  $\mathbf{c}$ . Our approach relies on results of [13], [28] developed for standard RIP, however, exploiting the block structure of  $\mathbf{c}$  leads to a much faster decay rate.

*Proposition 3:* Suppose  $\mathbf{D}$  is an  $n \times N$  matrix from the Gaussian ensemble, namely  $[\mathbf{D}]_{ik} \sim \mathcal{N}(0, \frac{1}{n})$ . Let  $\delta_{k|\mathcal{I}}$  be the smallest value satisfying the block RIP (21) over  $\mathcal{I} = \{d_1 = d, \dots, d_m = d\}$ , assuming  $N = md$  for some integer  $m$ . Then, for every  $\epsilon > 0$  the block RIP constant  $\delta_{k|\mathcal{I}}$  obeys (for  $n, N$  large enough, and fixed  $d$ )

$$\text{Prob} \left( \sqrt{1 + \delta_{k|\mathcal{I}}} > 1 + (1 + \epsilon)f(r) \right) \leq 2e^{-NH(r)\epsilon} \cdot e^{-m(d-1)H(r)}. \quad (67)$$

Here, the ratio  $r = kd/N$  is fixed,

$$f(r) = \sqrt{\frac{N}{n}} \left( \sqrt{r} + \sqrt{2H(r)} \right),$$

and  $H(q) = -q \log q - (1-q) \log(1-q)$  is the entropy function defined for  $0 < q < 1$ .

The assumption that  $d_i = d$  simplifies the calculations in the proof. Following the proof, we shortly address the more difficult case in which the blocks have varying lengths. We note that Proposition 3 reduces to the result of [13] when  $d = 1$ . However, since  $f(r)$  is independent of  $d$ , it follows that for  $d > 1$  and fixed problem dimensions  $n, N, r$ , block-RIP constants are smaller than the standard RIP constant. The second exponent in the right-hand side of (67) is responsible for this behavior.

*Proof of Proposition 3:* Let  $\lambda = (1 + \epsilon)f(r)$  and define

$$\bar{\sigma} = \max_{|T|=k,d} \sigma_{\max}(\mathbf{D}_T), \quad \underline{\sigma} = \min_{|T|=k,d} \sigma_{\min}(\mathbf{D}_T) \quad (68)$$

where  $\sigma_{\max}(\mathbf{D}_T)$ ,  $\sigma_{\min}(\mathbf{D}_T)$ , are the largest and the smallest singular values of  $\mathbf{D}_T$ , respectively. We use  $|T| = k, d$  to de-

note a column subset of  $\mathbf{D}$  consisting of  $k$  blocks of length  $d$ . For brevity, we omit subscripts and denote  $\delta = \delta_{k|\mathcal{I}}$ . The inequalities in the definition of block-RIP (21) imply that

$$1 + \delta \geq \bar{\sigma}^2 \tag{69}$$

$$1 - \delta \leq \underline{\sigma}^2. \tag{70}$$

Since  $\delta$  is the smallest number satisfying these inequalities we have that  $1 + \delta = \max(\bar{\sigma}^2, 2 - \underline{\sigma}^2)$ . Therefore

$$\begin{aligned} & \text{Prob}(\sqrt{1 + \delta} > 1 + \lambda) \\ &= \text{Prob}\left(\sqrt{\max(\bar{\sigma}^2, 2 - \underline{\sigma}^2)} > 1 + \lambda\right) \end{aligned} \tag{71}$$

$$\leq \text{Prob}(\bar{\sigma} > 1 + \lambda) + \text{Prob}(\sqrt{2 - \underline{\sigma}^2} > 1 + \lambda). \tag{72}$$

Noting that  $\underline{\sigma} \geq 1 - \lambda$  implies  $\sqrt{2 - \underline{\sigma}^2} \leq 1 + \lambda$  we conclude that

$$\begin{aligned} & \text{Prob}(\sqrt{1 + \delta} > 1 + \lambda) \\ & \leq \text{Prob}(\bar{\sigma} > 1 + \lambda) + \text{Prob}(\underline{\sigma} < 1 - \lambda). \end{aligned} \tag{73}$$

We now bound each term in the right-hand side of (73) using a result of Davidson and Szarek [45] regarding the concentration of the extreme singular values of a Gaussian matrix. It was proved in [45] that an  $m \times n$  matrix  $\mathbf{X}$  with  $n \geq m$  satisfies

$$\text{Prob}(\sigma_{\max}(\mathbf{X}) > 1 + \sqrt{m/n} + t) \leq e^{-nt^2/2} \tag{74}$$

$$\text{Prob}(\sigma_{\min}(\mathbf{X}) < 1 - \sqrt{m/n} - t) \leq e^{-nt^2/2}. \tag{75}$$

Applying a union bound leads to

$$\begin{aligned} & \text{Prob}\left(\bar{\sigma} > 1 + \sqrt{\frac{kd}{n}} + t\right) \\ & \leq \sum_{|T|=k,d} \text{Prob}\left(\sigma_{\max}(\mathbf{D}_T) > 1 + \sqrt{\frac{kd}{n}} + t\right) \end{aligned} \tag{76}$$

$$\leq \sum_{|T|=k,d} e^{-nt^2/2} \tag{77}$$

$$= \binom{m}{k} e^{-nt^2/2}. \tag{78}$$

Using the well-known bound on the binomial coefficient (for sufficiently large  $m$ )

$$\binom{m}{k} \leq e^{mH(k/m)} \tag{79}$$

we conclude that

$$\text{Prob}\left(\bar{\sigma} > 1 + \sqrt{\frac{kd}{n}} + t\right) \leq e^{mH(k/m)} e^{-nt^2/2}. \tag{80}$$

To utilize this result in (73) we rearrange

$$1 + \lambda = 1 + (1 + \epsilon)f(r) \tag{81}$$

$$= 1 + (1 + \epsilon)\left(\sqrt{\frac{kd}{n}} + \sqrt{\frac{2N}{n}H(r)}\right) \tag{82}$$

$$\geq 1 + \sqrt{\frac{kd}{n}} + \sqrt{(1 + \epsilon)\frac{2N}{n}H(r)} \tag{83}$$

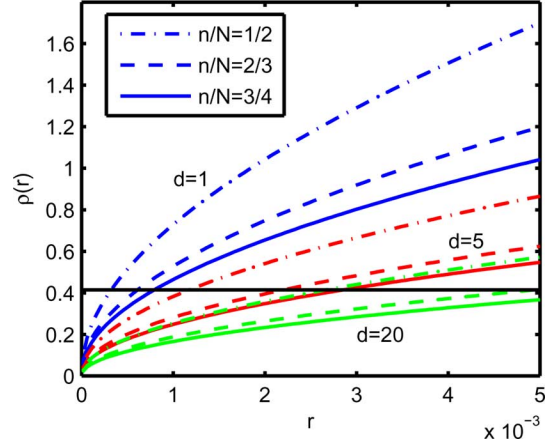


Fig. 4. The upper bound on  $\delta_{k|\mathcal{I}}$  as a function of the sparsity ratio  $r$ , for three sampling rates  $n/N$ , and three block structures  $d = 1, 5, 20$ . The horizontal threshold is fixed on  $\rho^* = \sqrt{2} - 1$  representing the threshold for equivalence derived in Theorem 1.

and obtain that

$$\begin{aligned} & \text{Prob}(\bar{\sigma} > 1 + \lambda) \\ & \leq \text{Prob}\left(\bar{\sigma} > 1 + \sqrt{\frac{kd}{n}} + \sqrt{(1 + \epsilon)\frac{2N}{n}H(r)}\right). \end{aligned} \tag{84}$$

Using (80) leads to

$$\text{Prob}(\bar{\sigma} > 1 + \lambda) \leq e^{mH(k/m)} e^{-\frac{n(1+\epsilon)2NH(r)}{2n}} \tag{85}$$

$$= e^{NH(r) - m(d-1)H(r) - (1+\epsilon)NH(r)} \tag{86}$$

$$\leq e^{-NH(r)\epsilon} e^{-m(d-1)H(r)}. \tag{87}$$

Similar arguments are used to bound the second term in (73), completing the proof.  $\square$

The proof of Proposition 3 can be adapted to the case in which  $d_i$  are not equal. In this case, the notation  $|T| = k, d$  is replaced by  $|T| = k|\mathcal{I}$  and has the following meaning:  $T$  indicates a column subset of  $\mathbf{D}$  consisting of  $k$  blocks from  $\mathcal{I}$ . Since  $\mathcal{I}$  contains variable-length blocks,  $|T|$  is not constant and depends on the particular column subset. Consequently, in order to apply the union bounds in (76) we need to consider the worst case scenario corresponding to the maximal block length in  $\mathcal{I}$ . Proposition 3 thus holds for  $d = \max(d_i)$ . However, it is clear that the resulting probability bound will not be as stringent as in the case of equal  $d_i = d$ , especially when the ratio  $\max(d_i)/\min(d_i)$  is large.

Proposition 3 holds as is for matrices  $\mathbf{D}$  from the Bernoulli ensemble, namely,  $[\mathbf{D}]_{ik} = \pm \frac{1}{\sqrt{n}}$  with equal probability. In fact, the proposition is true for any ensemble for which the concentration of extreme singular values holds.

The following corollary emphasizes the asymptotic behavior of block-RIP constants per given number of samples.

*Corollary 3:* Consider the setting of Proposition 3, and define  $g(r) = \sqrt{\frac{N}{n}}\left(\sqrt{r} + \sqrt{2H(r)d^{-1}}\right)$ . Then

$$\text{Prob}\left(\sqrt{1 + \delta_{k|\mathcal{I}}} > 1 + (1 + \epsilon)g(r)\right) \leq 2e^{-mH(r)\epsilon}. \tag{88}$$

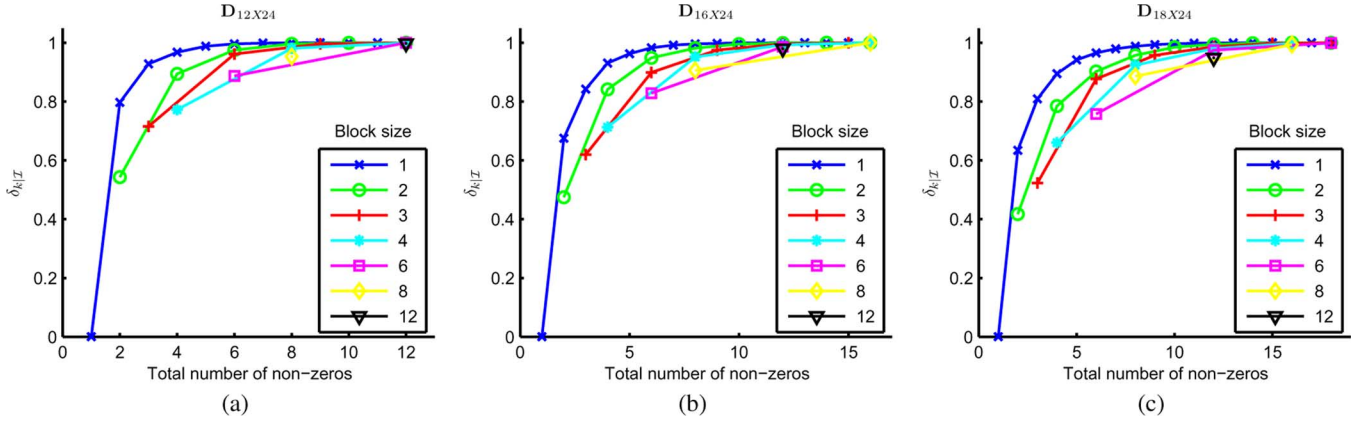


Fig. 5. The standard and block-RIP constants  $\delta_{k|\mathcal{I}}$  for three different dimensions  $n, N$ . Each graph represent an average over 10 instances of random matrix  $\mathbf{D}$ . Each instance of  $\mathbf{D}$  is scaled by a factor such that (18) is satisfied with  $\alpha + \beta = 2$ .

*Proof:* Let  $\lambda = (1 + \epsilon)g(r)$ . The result then follows by replacing (81)–(83) with

$$1 + \lambda \geq 1 + \sqrt{\frac{kd}{n}} + \sqrt{(1 + \epsilon)\frac{2N}{nd}H(r)}, \quad (89)$$

which leads to  $\text{Prob}(\bar{\sigma} > 1 + \lambda) \leq e^{-mH(r)\epsilon}$ .  $\square$

To evaluate the asymptotic behavior of block-RIP we note that for every  $\epsilon > 0$  the right-hand side of (88) goes to zero when  $N = md \rightarrow \infty$ . Consequently, for fixed  $d$

$$\delta_{k|\mathcal{I}} < \rho(r) \triangleq -1 + [1 + g(r)]^2 \quad (90)$$

with overwhelming probability. In Fig. 4, we compute  $\rho(r)$  for several problem dimensions and compare it with standard RIP which is obtained when  $d = 1$ . Evidently, as the nonzero entries are forced to block structure, a wider range of sparsity ratios  $r$  satisfy the condition of Theorem 1.

Although Fig. 4 shows advantage for block-RIP, the absolute sparsity ratios predicted by the theory are pessimistic as also noted in [13], [28] in the case of  $d = 1$ . To offer a more optimistic viewpoint, the RIP and block-RIP constants were computed brute-force for several instances of  $\mathbf{D}$  from the Gaussian ensemble. Fig. 5 plots the results and qualitatively affirms that block-RIP constants are more “likely” to be smaller than their standard RIP counterparts, even when the dimensions  $n, N$  are relatively small.

An important question is how many samples are needed roughly in order to guarantee stable recovery. This question is addressed in the following proposition, which quotes a result from [46] based on the proofs of [47]; we rephrase the result to match our notation.

*Proposition 4 ([46, Theorem 3.3]):* Consider the setting of Proposition 3, namely, a random Gaussian matrix  $\mathbf{D}$  of size  $n \times N$  and block sparse signals over  $\mathcal{I} = \{d_1 = d, \dots, d_m = d\}$ , where  $N = md$  for some integer  $m$ . Let  $t > 0$  and  $0 < \delta < 1$  be constant numbers. If

$$n \geq \frac{36}{7\delta} \left( \ln(2L) + kd \ln \left( \frac{12}{\delta} \right) + t \right) \quad (91)$$

where  $L = \binom{m}{k}$ , then  $\mathbf{D}$  satisfies the block-RIP (21) with restricted isometry constant  $\delta_{k|\mathcal{I}} = \delta$ , with probability at least  $1 - e^{-t}$ .

As observed in [46], the first term in (91) has the dominant impact on the required number of measurements in an asymptotic sense. This term quantifies the amount of measurements that are needed to code the exact subspace in which the sparse signal resides. Specifically, for block sparse signals

$$(m/k)^k \leq L = \binom{m}{k} \leq (em/k)^k. \quad (92)$$

Thus, for a given fraction of nonzeros  $r = kd/N$ , roughly  $n \approx k \log(m/k) = -k \log(r)$  measurements are needed. For comparison, to satisfy the standard RIP a larger number  $n \approx -kd \log(r)$  is required. Block sparsity reduces the total number of subspaces and therefore requires  $d$  times less measurements to code the signal subspace. The second term in (91) has a smaller contribution to the number of measurements. This term is proportional to  $kd$ , which is the number of nonzero values. Since the number of nonzeros is the same regardless of the sparsity structure, this term is not reduced in the block setting.

Note that Corollary 4 puts the emphasis on the required problem dimensions to satisfy a given RIP level. In contrast, Proposition 3 provides a tail bound on the expected isometry constant for given problem dimensions.

## VIII. CONCLUSION

In this paper, we studied the problem of recovering an unknown signal  $x$  in an arbitrary Hilbert space  $\mathcal{H}$ , from a given set of  $n$  linear samples. The signal  $x$  is known to lie in a union of subspaces, so that  $x \in \mathcal{V}_i$  where each of the subspaces  $\mathcal{V}_i$  is a sum of  $k$  subspaces  $\mathcal{A}_i$  chosen from an ensemble of  $m$  possibilities. While previous treatments of this model considered invertibility conditions, here we provide concrete recovery algorithms for a signal over a structured union.

We began by showing that recovering  $x$  can be reduced to a sparsity problem in which the goal is to recover a block-sparse vector  $\mathbf{c}$  from measurements  $\mathbf{y} = \mathbf{D}\mathbf{c}$  where the nonzero values in  $\mathbf{c}$  are grouped into blocks. The measurement matrix  $\mathbf{D}$  is equal

to  $S^*A$  where  $S^*$  is the sampling operator and  $A$  is a set transformation corresponding to a basis for the sum of all  $A_i$ . To determine  $\mathbf{c}$ , we suggested a mixed  $\ell_2/\ell_1$  convex optimization program that takes on the form of an SOCP. Relying on the notion of block-RIP, we developed sufficient conditions under which  $\mathbf{c}$  can be perfectly recovered using the proposed algorithm. We also proved that under the same conditions, the unknown  $\mathbf{c}$  can be stably approximated in the presence of noise. Furthermore, if  $\mathbf{c}$  is not exactly block-sparse, then its best block-sparse approximation can be approached using the proposed method. We then showed that when  $\mathbf{D}$  is chosen at random, the recovery conditions are satisfied with high probability.

Specializing the results to MMV systems, we proposed a new method for sampling in MMV problems. In this approach, each measurement vector depends on all the unknown vectors. As we showed, this can lead to better recovery rate. Furthermore, we established equivalence results for a class of MMV algorithms based on RIP.

Throughout the paper, we assumed a finite union of subspaces as well as finite dimension of the underlying spaces. An interesting future direction to explore is the extension of the ideas developed herein to the more challenging problem of recovering  $x$  in a possibly infinite union of subspaces, which are not necessarily finite-dimensional. Recovery methods for sparse signals in infinite dimensions have been addressed in some of our previous work [16]–[20]. In particular, we have shown that a signal lying in a union of shift-invariant subspaces can be recovered efficiently from certain sets of sampling functions. A first step in the direction of treating infinite unions of infinite subspaces is the example studied in [21] in which we treat an infinite union resulting from unknown time delays. In our future work, we intend to combine these results with those in the current paper in order to develop a more general theory for recovery from a union of subspaces.

A recent preprint [48] that was posted online after the submission of this paper proposes a new framework called model-based compressive sensing (MCS). The MCS approach assumes a vector signal model in which only certain predefined sparsity patterns may appear. In general, obtaining efficient recovery algorithms in such scenarios is difficult, unless further structure is imposed on the sparsity patterns. Therefore, the authors consider two types of sparse vectors: block sparsity as treated here, and a wavelet tree model. For these settings, they propose generalizations of several known greedy algorithms. The union model developed in this paper is broader than the block-sparse setting treated in [48] in the sense that it allows to model linear dependencies between the nonzero values rather than only between their locations, by appropriate choice of subspaces in (6), (7). In addition, we aim at optimization-based recovery algorithms (26), (47) which require selecting the objective in order to promote the model properties.

#### ACKNOWLEDGMENT

The authors would like to thank Volken Cevher for fruitful discussions regarding the MCS framework.

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