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# Nonconvex lp-Norm Regularized Sparse Self-Representation for Traffic Sensor Data Recovery

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**ABSTRACT** Recovering missing values from incomplete traffic sensor data is an important task for intelligent transportation system because most algorithms require data with complete entries as input. Selfrepresentation-based matrix completion attempts to optimally represent each sample by linearly combining other samples when conducting missing values recovery. Typically, it implements sparse or dense combination through imposing either *l*1-norm or *l*2-norm regularization over the representation coefficients, which is not always optimal in practice. To permit more flexibility, we propose in this paper a novel approach termed as *l*<sub>p</sub>-norm regularized sparse self-representation (SSR-*l<sub>p</sub>*) by incorporating nonconvex *l<sub>p</sub>*-norm with  $0 < p < 1$ as regularization. In such a way, it is able to produce more sparsity than  $l_1$ -norm and in turn facilitates the accurate recovery of missing data. We further develop an efficient iterative algorithm for solving SSR-*lp*. The performance of this method is evaluated on a real-world road network traffic flow data set. The experimental results verify the advantage of our method over other competing algorithms in recovering missing values.

**INDEX TERMS** Traffic sensor data, missing values, *lp*-norm regularization, sparse self-representation.

# **I. INTRODUCTION**

Traffic flow data collected by geographically distributed sensors has come to play an important role in advanced intelligent transportation system (ITS) because most traffic services provided by ITS depend on the accuracy and completeness of data. For example, short-term traffic flow forecasting [1], [2], which is of paramount importance for realizing proactive traffic control and effective route planning, requires data fed into specific predictive models, e.g. support vector machine [2], [3], neural networks [4], etc., to be complete without missing entries. Despite the fast growing reliability of sensing equipment and transmission network [5], [6], missing sensor data is still prevalent and inevitable in current ITS. For example, it was reported that for a dense road network in the city of Melbourne, about 8% of sensor can reach up to 56% missing data. Similarly, about 10% of daily traffic

flow in Beijing is missing. There are many reasons that lead to missing data, such as sensor malfunction, transmission distortion, and other unexpected exogenous factors. Without proper preprocessing, datum with missing entries cannot be directly utilized by most machine learning algorithms.

To address the above mentioned missing sensor data problem, many imputation methods have been proposed in the literature during the past decades. Here, imputation means the procedure that generates plausible estimations for the missing values (MVs) in a given incomplete data [7]. By means of MV imputation, the incomplete data can be converted into complete one and then used in traditional machine learning algorithms. Due to the connectivity of road network and the regularity of human travel activity, traffic sensor data collected at different time intervals and different road segments is essentially correlated with each other. Consequently, such

a kind of intrinsic correlation between traffic sensor data makes the recovery of missing values feasible and reliable in practice. Nowadays, some typical imputation methods have been developed in the literature, including K-nearest neighbors (KNN) [8], singularity value decomposition (SVD) [9], local least squares regression (LLS) [10], [11], probabilistic principle component analysis (PPCA) [12], [13], low-rank matrix completion (LRMC) [14]–[17], etc.

Recently, self-representation based matrix completion [18] was developed and shows competitive performance in comparison with other imputation methods. Essentially, selfrepresentation, as a general concept, refers to that each sample can be well represented as a linear combination of other samples with representation coefficients (weights) characterizing the contribution of other samples. Self-representation has already been widely exploited in some pattern recognition tasks, such as subspace clustering [19], [20], feature selection [21], etc., because of its simplicity and effectiveness. However, far less work exists on missing data recovery via self-representation. Different from self-representation with complete data, the representation of each sample cannot be obtained directly in the context of missing data. As a result, the recovery of missing data and the search for representation have to be implemented simultaneously. A key factor in selfrepresentation based matrix completion is the selection of suitable regularization on representation coefficients. After comparing  $l_2$ -norm,  $l_1$ -norm, and nuclear-norm [17], the previous work [18] verified from experiments that  $l_1$ -norm regularization performs much better than other forms of regularization in terms of recovery accuracy. The advantages of *l*1-norm regularization are two-fold. First, the representation vector of a target sample with respect to all other samples is sparse, implying that only a few of samples have nonzero coefficients. This is because  $l_1$ -norm is the tightest convex relaxation of  $l_0$ -norm [22]. Second,  $l_1$ -norm minimization generally leads to convex optimization problem [23] with efficient implementation. Nevertheless, a potential issue concerning *l*1-norm is that it may fail to find desired solution [22]. From this view of point,  $l_1$ -norm could be too restrictive and not sufficiently flexible, which in turn influences the imputation accuracy of missing data.

Inspired by the above discussion, in this paper, we propose a novel self-representation based matrix completion approach for missing data recovery by incorporating *lp*-norm regularization with  $0 < p < 1$  [24]. It has been observed that as a nonconvex surrogate of  $l_0$ -norm,  $l_p$ -norm minimization can often achieve more sparsity than *l*<sub>1</sub>-norm minimization [25], since it is closer to  $l_0$ -norm when  $p$  is smaller than 1. Theoretically,  $l_p$ -norm requires weaker conditions than  $l_1$ -norm to guarantee successful recovery of sparse signal [26]. We also take into account the nonnegative property of traffic sensor data through optimization, which is an inherent requirement for many real-world physical systems. The proposed method is termed as *lp*-norm regularized sparse self-representation (abbreviated as SSR-*l<sup>p</sup>* for brevity). Despite more flexibility with  $l_p$ -norm, it has difficulty in solving the resultant model because *lp*-norm is typically nonconvex. To address this issue, we further develop an efficient alternating algorithm which combines iterative reweighted least squares (IRLS) [22] as well as classic gradient projection (GP) method.

We summarize the main contributions of this paper as follows: (1) SSR-*l<sup>p</sup>* is a general framework benefiting from both self-representation and *lp*-norm regularization, providing flexible framework for MV imputation. (2) An efficient alternating optimization algorithm is proposed to solve SSR-*l<sup>p</sup>* model. (3) Extensive experiments on real-world traffic sensor data verify the effectiveness of our method in comparison with other related algorithms.

The rest of this paper is organized as follows. In Section II, we briefly review and analyze some popular methods for missing data recovery. In Section III, we present the proposed SSR-*l<sup>p</sup>* model and its optimization algorithm. Section IV reports the experimental results on real-world traffic sensor data. Finally, Section V gives the conclusions and discusses future work.

#### **II. RELATED WORKS**

So far, many MV imputation or recovery approaches have been proposed in the literature. These methods can be roughly classified into the following categories.

# A. REGRESSION BASED METHODS

The methods in this category attempt to characterize the relationship between missing values and observed values by regression models built based on training data. Regression models can be divided into parametric and nonparametric regression, thus further refining the division of imputation methods in this category. Some typical regression models include K-nearest neighbor (KNN) regression, least squares regression [10], support vector regression [27], neural networks [28], [29], [43], etc. For example, KNN imputation first select K nearest samples for the sample with MVs, followed which the MVs can be estimated as the weighted average of those selected samples. Following similar idea, local least squares (LLS) imputation [10], [30] also selects K nearest samples for the sample with MVs, but different from KNN imputation, it describes the relation between MVs and observed values by virtue of least squares regression, allowing more flexibility than simple weighted average. LLS has been proved to yield promissing performance in traffic sensor data recovery [31] and other domains [11].

### B. PROBABILISTIC MODEL BASED METHODS

In this type of methods, the complete data is supposed to follow a probabilistic distribution with specific form but unknown model parameters. Based on the observed values, both the model parameters and the missing data can be simultaneously estimated following maximum likelihood estimation (MLE) or full Bayesian framework [32], [33]. A popular algorithm to achieve such joint estimation is based on expectation-maximization (EM) [34]. A typical method belonging to this category is the so-called probabilistic

principal component analysis (PPCA) [13] assuming that the data follows multivariate Gaussian model. The missing data and model parameters are alternatively estimated by EM algorithm. This method has shown promising results in the imputation of traffic sensor data. This type of methods [13], [34], [35] imposes a global distribution assumption about data, thus being effective when data is consistent with the assumed distribution.

# C. MATRIX COMPLETION BASED METHODS

This type of methods organizes all samples into a matrix and achieve MVs recovery based on certain property of the matrix. One of the most well-known approach belonging to this class is low-rank matrix completion (LRMC) [14] which assumes the matrix is of low-rank structure. For traffic matrix, this assumption is reasonable to a certain extent, because traffic flows within the same road network are spatially and temporally correlated with each other. Owing to such inherent correlation, it was reported that LRMC is able to produce accurate recovery of missing data [36], [37], [44]. Recently, LRMC has attracted considerable attention and researchers have developed many optimization algorithms dedicated to solving LRMC model, such as SVT [14], FPCA [38], ADMM [39], etc. However, LRMC takes a global view on the data matrix, without sufficiently accounting for the difference between samples [17]. It may produce suboptimal recovery for samples with complex inherent structure, e.g., multiple subspaces. To this end, self-representation based matrix completion [18] was presented recently, aiming to characterize the relations between samples through linear combination. The missing data recovery and effective representation are solved jointly, indicating the two tasks can facilitate each other.

# **III. THE PROPOSED METHOD**

In this section, we first present the formulation of  $l_p$ -norm regularized sparse self-representation. Then, an efficient optimization algorithm for solving this model is proposed.

Formally, let  $X = [x_1, x_2, \dots, x_N] \in \mathbb{R}^{d \times N}$  be the given data matrix, where  $x_i \in \mathbb{R}^d$  denotes the *i*-th sample with *d* features, *N* is the total number of samples. Notice that in our problems, not all of the elements in *X* are observable. Due to diverse causes, many elements in *X* are missing. Here, we use  $\Omega$  to denote the indexes of missing values in *X*. Thereby, the central task of MV imputation is to estimate  $X_{\Omega}$ as accurately as possible.

# A. lp-NORM REGULARIZED SPARSE SELF-REPRESENTATION

In the spirit of self-representation [18], [19], each data sample can be represented as a linear combination of other samples. In particular, we have

$$
x_i \approx \sum_{j=1, j \neq i}^{N} w_i(j)x_j \tag{1}
$$

where  $w_i(j)$  denotes the combinatorial coefficient or weight of  $x_j$  in the resulting linear combination. Further, we introduce the following  $l_p$ -norm (0 $\lt p \lt 1$ ) regularization [22], [24] on the weight vector  $w_i = [w_i(1), w_i(2), \dots, w_i(N)]^T$ 

$$
||w_i||_p^p = \sum_{j=1}^N |w_i(j)|^p
$$
 (2)

with hope that most elements in  $w_i$  should be zero or close to zero such that the corresponding samples can be eliminated from the representation of *x<sup>i</sup>* .

Besides the above task, another difficulty comes from the fact that many elements in *X*, i.e.  $X_{\Omega}$ , are unknown while only the rest part of *X* are observed. As a result, it is infeasible to find weight vector  $w_i$  directly given a set of incomplete data samples. In other words, we need to discover the above sparse linear representation structure among data and meanwhile estimate the involved missing values. In fact, it is expected that the reliable estimation of MVs and the discovery of sparse linear relations among data are related and thus would benefit from each other, indicating the two tasks can be solved in a uniformed framework. In addition, for many applications, the data that real physical system records is usually nonnegative [6]. As a result, nonnegativity should also be taken into account. Based on the above discussion, we present *lp*-norm regularized sparse self-representation (SSR-*lp*) for MV imputation as follows

<span id="page-2-0"></span>
$$
\min_{X_{\Omega}, W} \frac{1}{2} \sum_{i=1}^{N} \|x_i - \sum_{j=1, j \neq i}^{N} w_i(j)x_j\|^2 + \lambda \sum_{i=1}^{N} \|w_i\|_p^p
$$
\n
$$
\text{s.t.} \quad X_{\Omega} \ge 0 \tag{3}
$$

where  $\lambda > 0$  is a prameter controlling the strength of  $l_p$ -norm regularization.

Let  $W = [w_1, w_2, \dots, w_N]$  and diag(*W*) stand for the diagonal elements of *W*. The above problem [\(3\)](#page-2-0) can be expressed in matrix form as follows

<span id="page-2-1"></span>
$$
\min_{X_{\Omega}, W} \frac{1}{2} \|X - XW\|^2 + \lambda \sum_{i=1}^N \|w_i\|_p^p
$$
  
s.t.  $X_{\Omega} \ge 0$ , diag(W) = 0 (4)

Notice that in model [\(4\)](#page-2-1), both  $X_{\Omega}$  and *W* are the decision variables, differentiating it from traditional subspace clustering [19], [20] where only *W* is the variable need to solve. It should be pointed out that problem [\(4\)](#page-2-1) naturally reduces to the models developed in [18] given  $p = 1$  or  $p = 2$ .

# B. OPTIMIZATION ALGORITHM

The problem [\(4\)](#page-2-1) does not allow a closed-form solution because of the coupling between decision variables  $X_{\Omega}$  and *W*, which leads the problem difficult to solve. However, we observe that the problem can be simplified if only one variable is optimized each time while fixing the other one. To this end, we develop an iterative algorithm to solve [\(4\)](#page-2-1) by alternatively optimizing over  $X_{\Omega}$  and *W* while holding the other variable fix.

# **Algorithm 1** Solve  $X_{\Omega}$  When *W* Is Fixed



- 1: Initialize  $X_{\Omega}$ , let  $c = 10^{-4}$
- 2: **while** not converged **do**
- 3: Compute gradient  $\nabla g(X_{\Omega})$  of  $g(X_{\Omega})$
- 4: Find step-size *l* with Armijo rule, i.e., choose  $l = \min\{1, \frac{1}{2}, \frac{1}{2^2}\}$  $\frac{1}{2^2}, \ldots$ } such that  $g(X_{\Omega}^{*}) \leq g(X_{\Omega}) + c \cdot trace((X_{\Omega}^{*} - X_{\Omega})^{T} \nabla g(X_{\Omega}))$ where  $X_{\Omega}^{*} = \max\{X_{\Omega} - l\nabla g(\hat{X}_{\Omega}), 0\}$ 5: Update MV estimation as  $X_{\Omega} \leftarrow X_{\Omega}^*$
- 

# 6: **end while**



Concretely, we first fix *W* and seek for the optimal solution of  $X_{\Omega}$ . In such a case, the problem [\(4\)](#page-2-1) is equivalent to the following constrained problem

<span id="page-3-0"></span>
$$
\min_{X_{\Omega}} \frac{1}{2} \|X - XW\|^2
$$
  
s.t.  $X_{\Omega} \ge 0$  (5)

In order to solve [\(5\)](#page-3-0), we develop an iterative algorithm based on gradient projection [23]. Let the derivative of  $g(X_{\Omega}) = \frac{1}{2} \|X - \overline{XW}\|^2$  with respect to  $X_{\Omega}$  be denoted by  $\nabla g(X_{\Omega}) = (X(I - W)(I - W)^{T})_{\Omega}$ , then the algorithm for solving [\(5\)](#page-3-0) can be summarized in **Algorithm 1**.

Secondly, we attempt to optimize *W* while holding  $X_{\Omega}$ . In such a case, the original problem [\(4\)](#page-2-1) becomes

<span id="page-3-1"></span>
$$
\min_{W} \frac{1}{2} \|X - XW\|^2 + \lambda \sum_{i=1}^{N} \|w_i\|_p^p
$$
  
s.t. diag(W) = 0 (6)

Notice that problem [\(6\)](#page-3-1) is separable with respect to the columns of *W*, therefore, we can solve each  $w_i$  by

<span id="page-3-2"></span>
$$
\min_{w_i} \frac{1}{2} \|x_i - Xw_i\|^2 + \lambda \|w_i\|_p^p
$$
  
s.t.  $w_i(i) = 0$  (7)

In order to handle the constraint in [\(7\)](#page-3-2), we denote  $\overline{X}_i = [x_1, \ldots, x_{i-1}, x_{i+1}, \ldots, x_N] \in \mathbb{R}^{d \times (N-1)}$ ,  $\overline{w}_i =$  $[w_i(1), \ldots, w_i(i-1), w_i(i+1), \ldots, w_i(N)]^T$ , then, the constrained problem [\(7\)](#page-3-2) can be converted into the unconstrained problem as follows

<span id="page-3-3"></span>
$$
\min_{\overline{w}_i} \frac{1}{2} \|x_i - \overline{X}_i \overline{w}_i\|^2 + \lambda \|\overline{w}_i\|_p^p \tag{8}
$$

As we can see, a major difficulty is that problem [\(8\)](#page-3-3) is nonconvex when  $0 < p < 1$ . In this work, in order to take the advantage of the problem structure, we present an algorithm targeting at [\(8\)](#page-3-3) by means of iteratively reweighted least squares (IRLS) [22] which is a popular technique in optimization. Specifically, by computing the derivative of the objective [\(8\)](#page-3-3) with respect to  $\overline{w}_i$  and setting to zero, we have

$$
\overline{X}_i^T(\overline{X}_i\overline{w}_i - x_i) + \frac{\lambda p \overline{w}_i(j)}{(\overline{w}_i(j)^2 + \epsilon)^{1-\frac{p}{2}}} = 0
$$
 (9)

where  $\epsilon$  is a small number to avoid division by zeros.



- 1: **while** not converged **do**
- 2: Use **Algorithm 1** to update current MVs  $X_{\Omega}$
- 3: **for**  $i = 1, 2, ..., N$  **do**
- 4: Compute  $D_t$  by using [\(11\)](#page-3-4)<br>5: Update  $w_i$  by using (10) or
- Update  $w_i$  by using [\(10\)](#page-3-5) or [\(13\)](#page-3-6)
- 6: **end for**
- 7: **end while**
- 8: **Output** the estimated missing data  $X_{\Omega}$

Then, the iterative procedure for solving  $\overline{w}_i$  is given by

<span id="page-3-5"></span>
$$
\overline{w}_i = (\overline{X}_i^T \overline{X}_i + \lambda D_t)^{-1} \overline{X}_i^T x_i
$$
\n(10)

where  $D_t$  is a diagonal matrix defined as

<span id="page-3-4"></span>
$$
D_{t} = \text{diag}(\frac{p}{(\overline{w}_{i}^{t}(j)^{2} + \epsilon)^{1-\frac{p}{2}}}),
$$
  
  $j = 1, ..., i - 1, i + 1, ..., N$  (11)

and  $\overline{w}_i^t$  is the solution at the *t*-th iteration.

Note that when the total number of samples is large, e.g.  $N \gg d$ , the inversion of matrix  $\overline{X}_i^T \overline{X}_i + \lambda D_t$  in (10) is computationally expensive, i.e.,  $O(N^3)$  in time complexity. To address this problem and make our algorithm feasible given large number of samples, we apply the well-known Sherman-Morrison-Woodbury formula [40] shown below

<span id="page-3-7"></span>
$$
(UCV + A)^{-1} = A^{-1} - A^{-1}U(C^{-1} + VA^{-1}U)^{-1}VA^{-1}
$$
 (12)

By combing [\(10\)](#page-3-5) and [\(12\)](#page-3-7), the optimal solution of  $\overline{w}_i$  can be rewritten as

<span id="page-3-6"></span>
$$
\overline{w}_i = \frac{1}{\lambda} D_t^{-1} - \frac{1}{\lambda} D_t^{-1} \overline{X}_i^T (\lambda I + \overline{X}_i D_t^{-1} \overline{X}_i^T)^{-1} \overline{X}_i D_t^{-1}
$$
(13)

where *I* is a  $d \times d$  identity matrix. In formula [\(13\)](#page-3-6), we only need to solve the inversion of a matrix with size  $d \times d$ . To this end, when  $N \gg d$ , the time complexity for solving  $w_i$ can be significantly reduced from  $O(N^3)$  to  $O(d^3)$ . The time complexity for solving all of  $w_i$  can be estimated as  $O(Nd^3)$ , which is linear with respect to the number of samples.

Finally, the whole iterative algorithm for solving SSR-*l<sup>p</sup>* model [\(4\)](#page-2-1) is summarized in **Algorithm 2**.

# **IV. EXPERIMENTS AND ANALYSIS**

# A. DATA DESCRIPTION

In this study, we evaluate the proposed SSR-*l<sup>p</sup>* algorithm on a real-world traffic flow dataset. The data was collected from Interstate 205 (I205) highways, serving the Portland-Vancouver metropolitan area in Oregon and Washington states, USA. The selected sub-area road network is shown in Fig. [1.](#page-4-0) Thirty inductive loop detectors which records the vehicle volume counts are chosen. The aggregation period is 15 minutes, thus yielding 96 sampling points in each day. In other words, each data sample can be viewed a point in a 96 dimensional space. The collection time period used in this study was from Mar. 1st to Aug. 31st in



<span id="page-4-0"></span>**FIGURE 1.** The selected sub-area road network of Portland, OR, USA.



<span id="page-4-1"></span>**FIGURE 2.** Illustration of traffic flow profiles from 30 detectors in the same day.

the year 2015. The data is publicly available at website (http://portal.its.pdx.edu/). After excluding weekends as well as holidays, we finally get volume data of 97 days. Finally, the total number of volume counts reaches  $96 \times 30 \times 97 =$ 279360. The whole traffic sensor data is organized as a 96  $\times$ 2910 matrix with each column representing a data sample. In Fig. [2,](#page-4-1) we illustrate 30 data samples, each of which was captured by a distinct detector in the same day. Note that the horizontal axis and the vertical axis in Fig. [2](#page-4-1) represent time and traffic volume, respectively. These samples intuitively reflect the traffic flow profiles at different road segments. As can be seen, despite of overall similarity among traffic flow profiles, it does exhibit some distinctions with respect to the variation patterns of different detectors. For instance, the maximum flow of some detectors is significantly larger than that of other detectors. In addition, the traffic flow at certain detectors clearly shows two peaks at rush hours while it is not very notable for other detectors. These slight yet important differences pose great challenge for MV imputation problem and render us to develop more flexible model such that the homogeneity as well as heterogeneity can be stimutanlously taken into accout.

#### B. CONFIGURATION

To comprehensively compare different methods, beside the proposed SSR-*lp*, we also include some closely-related



<span id="page-5-1"></span><span id="page-5-0"></span>**FIGURE 3.** Convergence curve under different missing patterns. (a) MAR. (b) MIXED.

TABLE 1. Imputation error (×10<sup>-2</sup>) under MCAR missing pattern.

Method		<b>KNN</b>	<b>LLS</b>	<b>PPCA</b>	<b>LRMC</b>	$SR-l2$	<b>SRSp</b>	$SSR-l_n$
0.1	Mean	97.40	73.40	81.28	86.15	81.35	55.72	$57.03(p=0.8)$
	Std	0.85	0.87	0.73	0.68	0.75	0.49	0.60
0.2	Mean	109.34	79.92	85.50	88.52	83.87	61.21	$60.91(p=0.8)$
	Std	0.56	0.39	0.38	0.37	0.41	1.12	0.55
0.3	Mean	121.96	87.89	88.11	91.05	86.70	71.99	$65.21(p=0.2)$
	Std	0.58	0.45	0.23	0.22	0.21	0.29	0.42
0.4	Mean	134.50	97.46	91.75	94.12	89.93	78.99	$71.79(p=0.2)$
	Std	0.12	0.39	0.22	0.27	0.21	0.43	0.18
0.5	Mean	147.16	110.86	96.04	97.82	93.69	87.56	$81.44(p=0.2)$
	Std	0.22	0.77	0.25	0.21	0.25	0.52	0.41

algorithms, including KNN, LLS [10], PPCA [13], LRMC [41], SR-*l*<sup>2</sup> [18] and SRSp [18]. Note that SRSp can be viewed as a special case of SSR-*l<sup>p</sup>* when *p* equals to 1. These methods covers the mainstream techniques for traffic data imputation, such as regression model, probabilistic model, etc. All of these algorithms were implemented in MATLAB environment on a PC with Intel(R) Core(TM) i7-4712MQ CPU and 12GB DDR4 RAM. There are some parameters involved in each method, such as the number of nearest neighbors for KNN and LLS, the subspace dimensionality for PPCA, etc. Following previous stuides [18], we adjust the parameters in each method such that best performance is achieved.

In order to simulate MVs and evaluate the imputation performance, we artificially generate missing entries for the data. Specifically, three missing patterns [13], [17] are considered in the experiment. (i) missing completely at random (MCAR) where the data points to be missing are completely independent of each other and occur as a set of isolated points randomly distributed, (ii) missing at random (MAR) where the appearance of missing points depends on their neighboring points. Therefore, this type of missing pattern looks like a group of successive MVs, (iii) a mixture of MCAR and MAR (MIXED), where half of MVs obey MCAR and the other half are from MAR.

We measure the recovery performance of each method by root mean squared error (RMSE) between the estimated values and the real values for those missing entries. Clearly, smaller RMSE indicates better recovery performance. We also define the missiong ratio  $\delta$  as the ratio of

the number of missing entries to the total number of entries. Moreover,  $\delta$  is changed from 0.1 to 0.5 with step 0.1 in order to investigate the variation of recovery performance against different missing ratios.

# C. CONVERGENCE ANALYSIS

In this work, an iterative algorithm alternatively recovering the missing data and optimizing sparse representation coefficients is developed to solve the proposed SSR-*l<sup>p</sup>* model. Next, we empirically investigate the convergence behavior of this algorithm under varying missing ratios and different missing patterns. Some convergence curves obtained in the experiments are shown in Fig[.3](#page-5-0) where the x-axis denotes the number of iterations and the y-axis denotes the logarithm of objective function. From these results, we can observe that our algorithm reduces the objective [\(13\)](#page-3-6) in each iteration, regardless of specific missing ratio and missing pattern. Moreover, the iterative algorithm we develop is able to converge quickly, usually requiring about 10-40 iterations in most cases.

# D. IMPUTATION PERFORMANCE COMPARISON

Considering the randomness when artificially simulating missing entries, we repeat each experiment five times and calculate the average imputation error (Mean) as well as the associated standard deviation (Std). The experimental results under MCAR, MAR, and MIXED missing patterns are reported in Tables [1,](#page-5-1) [2](#page-6-0) and [3,](#page-6-1) respectively. Note that the number in parenthesis of SSR-*l<sup>p</sup>* column indicates the value





<span id="page-6-2"></span><span id="page-6-0"></span>**FIGURE 4.** Imputation results obtained by KNN.

	<b>TABLE 2.</b> Imputation error $(x10^{-2})$ under MAR missing pattern.					
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Method		<b>KNN</b>	<b>LLS</b>	<b>PPCA</b>	<b>LRMC</b>	$SR-l_2$	<b>SRSp</b>	$SSR-l_n$
0.1	Mean	127.69	82.84	92.87	105.28	98.27	65.62	$66.20(p=0.8)$
	Std	1.30	0.58	1.03	1.19	1.09	0.91	0.49
0.2	Mean	135.73	90.40	96.21	105.37	97.97	74.08	$71.03(p=0.8)$
	Std	1.98	1.05	0.98	0.93	1.00	1.60	0.99
0.3	Mean	143.45	98.82	98.75	106.70	99.11	85.67	$78.36(p=0.6)$
	Std	0.79	0.79	0.41	0.55	0.40	1.54	0.78
0.4	Mean	150.82	110.23	101.66	108.44	101.14	93.38	$85.54(p=0.6)$
	Std	0.66	1.12	0.37	0.40	0.45	0.82	0.72
0.5	Mean	158.09	126.44	105.00	110.79	104.14	100.37	$94.22(p=0.2)$
	Std	0.38	0.88	0.57	0.41	0.54	0.54	0.83

<span id="page-6-1"></span>TABLE 3. Imputation error (×10<sup>-2</sup>) under MIXED missing pattern.



of *p* used in this experiment. As we can see, overall, MCAR missing pattern is the easiest situation while MAR is the most difficult case in terms of imputation. It is reasonable because successive missing will lose much valuable information about correlation, thus increasing the difficulty of accurate recovery. With respect to recovery performance, we find that KNN performs worst among these algorithms. LLS, PPCA, LRMC, SR-*l*2, and SRSp all significantly outperform KNN. In particular, LLS works well in low missing ratio, however, rapidly degrades when missing ratio increases. Self-representation based methods, including SR-*l*2, SRSp, and our proposed SSR-*lp*, obtain superior imputation performance than other competing methods. Comparing these three methods, we find that SRSp obtains better performance than SR-*l*2, indicating that sparsity is a crucial factor for self-representation based imputation. This conclusion is consistent with that drawn in [18]. At last, SSR-*l<sup>p</sup>* achieves best performance in most cases. In fact, it is interesting to notice that as the missing ratio increases, a smaller *p* is preferred which implies fewer samples should be selected for MV recovery in such situations. Some examples under MIXED missiont pattern and  $\delta = 0.3$  are shown in Fig[.4-](#page-6-2)Fig[.10.](#page-8-0) As we can see, the proposed SSR-*l<sup>p</sup>* achieves small residual in the recovery of missing data.

# E. INFLUENCE OF  $I<sub>D</sub>$  ON PERFORMANCE

In what follows, we investigate the recovery performance when varying the value of  $p$  in order to confirm that it is an



**FIGURE 5.** Imputation results obtained by LLS.



**FIGURE 6.** Imputation results obtained by PPCA.



**FIGURE 7.** Imputation results obtained by LRMC.

important factor for self-representation based matrix completion. In particular, we change the parameter  $p$  in the range of {0.2, 0.4, 0.6, 0.8, 1.0} and record the best performance

for each candidate value. Some experimental results under MCAR, MAR and MIXED missing patterns are presented in Fig[.11.](#page-9-0) As we can see, the variation trends of performance









**FIGURE 8.** Imputation results obtained by SR-/<sub>2</sub>.



**FIGURE 9.** Imputation results obtained by SRSp.



<span id="page-8-0"></span>**FIGURE 10. Imputation results obtained by SSR-** $l_p$ **.** 

when varying *p* is different under different missing ratio δ. When  $\delta$  is small, the performance can be improved by increasing the value of  $p$ . In contrast, when  $\delta$  is large, smaller  $p$ 



800

700

600

500

 $\overline{0}$ 

 $\overline{\mathbf{0}}$ 

100

200

300

400

Observed value  $(b)$ 

500

600

700

800



is preferred. The results indicate that given more observed entries, our method is apt to use more samples for accurate imputation. On the contrary, when there are too many missing



FIGURE 11. Imputation performance variation with respect to different value of the parameter *p*. (a) MCAR,  $δ = 0.1$ . (b) MCAR,  $δ = 0.3$ . (c) MCAR,  $\delta = 0.5$ .

<span id="page-9-0"></span>

**FIGURE 12.** Variation of weights when increasing p from 0.2 (top row) to 1.0 (bottom row) with step 0.2.

values, our method adjusts itself to utilize samples as fewer as possible. These results are essentially consistent with the above quantitative comparison.

Next, we investigate how the parameter *p* influences the sparsity of the resulting model. In this experiments, we take one data sample as instance and fix  $\lambda$  to a constant and change *p* in the same range above. Fig. 6 demonstrates the resulting  $w_1 \in \mathbb{R}^{2910}$  after optimization. We show the weights with absolute value larger than  $10^{-4}$  following [42], [22]. As we can see, when *p* equals to 0.2, only four samples are selected. When *p* increases to 0.4, three extra samples are selected. In a similar way, with the increase of *p*, more and more samples are selected. These results empirically verify that, the proposed  $SSR-l_p$  model is able to produce the solution with more sparsity given a small value of *p*.

# **V. CONCLUSIONS**

In this paper, we develop a novel MV imputation algorithm based on self-representation and *lp*-norm minimization. With the introduction of  $l_p$ -norm, our method is able to find sparser representation for each sample, which in turn facilitate the accurate recovery of missing data. To solve the resulting model, we further develop an algorithm which optimizes the missing data and the representation coefficients alternatively.

The experimental results confirm the effectiveness of our method. An interesting extension of our work is investigating the nonlinear formulation of the proposed model, which is more powerful in modeling nonlinear structure of data. Another future work is to extend our proposal to large-scale problems by applying some techniques, such as parallelization, etc.

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