

Provable Low-Rank Plus Sparse Matrix Recovery Via Nonconvex Regularizers

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Abstract

This paper considers a large class of problems where we seek to recover a low rank matrix and/or sparse vector from some set of measurements. While methods based off of convex relaxations suffer from a (possibly large) estimator bias, and other nonconvex methods require the rank or sparsity to be known a priori, we utilize nonconvex regularizers to minimize the rank and l_0 norm without the estimator bias from the convex relaxation. We present a novel analysis of the alternating proximal gradient descent algorithm applied to such problems, and bound the error between the iterates and the ground truth sparse and low rank matrices. The algorithm and error bound can be applied to sparse optimization, matrix completion, and robust principal component analysis as special cases of our results.

1. Introduction

In order to better understand large data-set and to make inferences about them, it is helpful to understand the underlying patterns in the dataset. Even when the underlying pattern is highly nonlinear, the data matrix can be approximated as being low rank, an observation that enables techniques to analyze the data in terms of a low dimensional latent space, such as principal Component Analysis (PCA), identifying outliers through Robust PCA (RPCA), and accurately inferring data points from very few observations of a data matrix through matrix completion.

Data analysis techniques based upon this low rank property have received much attention in the past decade, with impressive computational results on large matrices and theoretical results guaranteeing the success of RPCA and matrix completion. Many of these results are based off of minimizing the nuclear norm of a matrix (defined as the sum of the singular values) as a surrogate for the rank function,

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similar to minimizing the l_1 norm to promote sparsity in a vector.

While the convex relaxation is an incredibly useful technique in many applications, minimizing the nuclear norm of a matrix has been shown to introduce a (sometimes very large) estimator bias. Intuitively, we expect to see this bias because if we hope to recover a rank r matrix, we must impose enough weight on the nuclear norm term so that the $(r + 1)$ th singular value is zero. By the nature of the nuclear norm, this requires also putting weight on minimizing the first r singular values, resulting in a bias towards zero proportional to the spectral norm of the noise added to the true data matrix.

Fortunately, recent work has shown that the estimator bias from convex regularizers can be reduced (or even eliminated, for well conditioned matrices) by using nonconvex regularizers, such as the Schatten- p norm or the minimax concave penalty (MCP). It has been shown that for sparse optimization, the nonconvexity introduced from these regularizers does not create a further burden in the the optimization process – in the right circumstances, the nonconvex problem has just one minimizer. Similar results for rank minimization problems have been previously unavailable, a gap that we have aimed to fill in this paper.

1.1. Summary of Contributions

In this paper, we focus on the nonconvex, unconstrained optimization problem where we find a low-rank matrix $L \in \mathbb{R}^{d_1 \times d_2}$ and sparse vector $s \in \mathbb{R}^{d_2}$.

$$\min_{L,s} \frac{\lambda_L}{d_1 d_2} \Phi_{\gamma_L}(L) + \frac{\lambda_s}{d_s} \phi_{\gamma_s}(s) + \frac{1}{2n} \|A_L(L) + A_S s - b\|_2^2 \quad (1)$$

We denote $\phi : \mathbb{R} \rightarrow \mathbb{R}_+$ to be a concave function used to promote sparsity in both the singular values of L and individual entries in s . We overload the notation to allow for ϕ_{γ_s} to be a function of a vector $x \in \mathbb{R}^{d_s}$ whose range is \mathbb{R}_+ , and we denote $\Phi_{\gamma_L} : \mathbb{R}^{d_1 \times d_2} \rightarrow \mathbb{R}_+$ as a surrogate to the rank function:

$$\phi_{\gamma_s}(x) = \sum_{i=1}^{d_s} \phi_{\gamma_s}(x_i), \quad \Phi_{\gamma_L}(X) = \sum_{i=1}^{\min(d_1, d_2)} \phi_{\gamma_L}(\sigma_i(X)).$$

where $\sigma_i(X)$ denotes the i th largest singular value of X . We restrict our focus to nonconvex regularizers that

are *amenable regularizers*, as described in (Loh & Wainwright, 2015), and defined below. Some key properties of amenable regularizers are stated in Appendix D.

Definition 1. A function $\phi_\gamma(t) : \mathbb{R}_+ \rightarrow \mathbb{R}_+$ is *amenable* if it satisfies the following criteria.

1. $\phi_\gamma(0) = 0$
2. ϕ_γ is non decreasing
3. For $t > 0$, the function $\frac{\phi_\gamma(t)}{t}$ is non increasing in t .
4. the function ϕ_γ is differentiable for all $t \neq 0$ and sub-differentiable at $t = 0$ with $\lim_{t \rightarrow 0^+} \phi'_\gamma(t) = 1$.
5. The function $\phi_\gamma(t)$ is ν weakly convex. That is, the function $\rho_\nu := \phi_\gamma(t) + \frac{\nu}{2}t^2$ is convex.

The linear mappings $\mathcal{A}_L : \mathbb{R}^{d_1 \times d_2} \rightarrow \mathbb{R}^n$ and $A_s \in \mathbb{R}^{d_s \times n}$ serve as the *observation models* of the underlying low rank matrices and sparse vectors. Most commonly, we are interested in the observation model

$$[\mathcal{A}_{\Omega^{obs}}(X)]_k = X_{i_k, j_k}$$

for $(i_k, j_k) \in \Omega^{obs}$, where $\Omega^{obs} \subseteq \{1, \dots, d_1\} \times \{1, \dots, d_2\}$ is the set of indices where we have a measurement of the low rank matrix we hope to reconstruct.

We present a very simple alternating algorithm to find a stationary point of (1). Though similar algorithms have been previously studied and shown to converge, we present a novel analysis of this algorithm and show that not only does this algorithm linearly converge, but it converges to the exact low rank matrix L^* and sparse vector s^* when no Gaussian noise is present. Furthermore, when Gaussian noise is present, we obtain an error bound that matches the minmax optimal rate. Our bound greatly improves upon bounds obtained in previous results analysing the convex relaxation and quantifying the observation based on computational results that nonconvex regularizers reduce the impact of noise on the quality of the estimator.

1.2. Related Works

The matrix completion problem is as follows: given the values of a low rank matrix for only a sparse set of indices, we seek to determine the rest of the values of the matrix. While the problem of finding the minimum rank matrix that fits the observations is NP hard in general, it has been shown that under some assumptions, the global minimizer to the convex problem

$$\min_{X \in \mathbb{R}^{d_1 \times d_2}} \|X\|_* \text{ s.t. } X_{ij} = M_{ij} \forall (i, j) \in \Omega^{obs}$$

is exactly M , where Ω^{obs} is the set of indices of M we have observed. If M is rank r and μ -incoherent (as defined in section 3.1), then with high probability for a set, Ω^{obs} of n indices chosen uniformly at random, M is the unique minimizer to the convex relaxation so long

as $n > C\mu r d_1 \log^{1.2}(d_1)$ for some universal constant C (Candes & Tao, 2010; Candès & Recht, 2012). This condition was later improved to $n > C\mu r d_1 \log(d_1)$ (Recht, 2011).

Using a convex relaxation for Robust PCA has similar results. While PCA is a powerful technique, it has been shown to be less reliable when just a sparse set of data points are grossly corrupted, and so the goal of RPCA is to identify and remove such corruptions by separating the data matrix into the sum of a low rank and sparse matrix.

$$\min_{L, s} \|L\|_* + \lambda_0 \|S\|_1 \text{ s.t. } L_{ij} + S_{ij} = M_{ij} \forall (i, j) \in \Omega^{obs}$$

The convex relaxation was shown to give the exact solution when every entry of M is observed in (Chandrasekaran et al., 2011), and when only partially observed (under the same assumptions necessary in matrix completion) by (Candès et al., 2011).

In many cases of practical interest, our measurements may have some level of noise in addition to being only partially observed or having some corrupted entries. For matrix completion, we can relax the constraint using a penalty formulation as follows:

$$\min_{X \in \mathbb{R}^{d_1 \times d_2}} \lambda \|X\|_* + \sum_{(i, j) \in \Omega^{obs}} (X_{ij} - M_{ij})^2$$

Likewise, RPCA can be formulated as solving:

$$\begin{aligned} \min_{L, S \in \mathbb{R}^{d_1 \times d_2}} \lambda_L \|L\|_* + \lambda_s \|S\|_1 \\ + \sum_{(i, j) \in \Omega^{obs}} (L_{ij} + S_{ij} - M_{ij})^2 \end{aligned}$$

Statistical guarantees on the performance of the first of these estimators are discussed in (Candes & Plan, 2010; Negahban & Wainwright, 2012; 2011), and the later in (Agarwal et al., 2012). Specific bounds and assumptions for these works are discussed in Section 4.

In order to reduce the estimator bias for l_1 minimization, (Candes et al., 2008) proposed an iteratively reweighted l_1 norm method to place more weight on minimizing smaller entries, and less on entries further from zero. This idea was generalized to minimizing any Amenable regularizers to promote sparsity. Theoretical results on the subject include algorithmic guarantees similar to the ones presented in this paper (Wang et al., 2014), and a proof that the non-convex problem has no spurious local minimizers (Loh & Wainwright, 2015; 2017).

The same regularizers could be used as a surrogate to the rank function, as originally proposed by (Mohan & Fazel, 2012). In (Lu et al., 2016; Canyi Lu, 2015), the authors propose a generalization of the singular value thresholding algorithm proposed by (Cai et al., 2008), which was

Table 1. Table of common nonconvex regularizers used for l_0 and rank minimization, along with the associated proximal operator.

	$\phi_\gamma(x)$	$\text{prox}_{\phi_\gamma}^\tau(y)$
l_1	$ x $	$\text{sign}(y)(y - \frac{1}{\tau})_+$
MCP	$\begin{cases} x - \frac{x^2}{2\gamma} & x \leq \gamma \\ \frac{\gamma}{2} & x > \gamma \end{cases}$	$\text{sign}(y) \min\left(y , \frac{\gamma\tau}{\gamma\tau-1}(y - \frac{1}{\tau})_+\right)$

later applied to the problem of RPCA in (Chartrand, 2012; Kang et al., 2015). For the problem of matrix completion, the algorithms proposed by (Yao et al., 2019) and (Sagan & Mitchell, 2020) achieve the fastest computational complexity in other state of the art methods.

Other approaches to low rank optimization rely on, instead of minimizing a surrogate to the rank function, constraining the matrix to be a given rank. This can be done using constrained optimization to optimize over the set of all rank r matrices (Vandereycken, 2013), or by using the low-rank factorization of a matrix $X = UV^T$ for $U \in \mathbb{R}^{d_1 \times r}$ and $V \in \mathbb{R}^{d_2 \times r}$ (Wen et al., 2012; Shen et al., 2014).

Previous work on theoretical results pertaining to rank and sparsity constrained methods have consisted of algorithmic guarantees ensuring that we can obtain a matrix sufficiently close to the ground truth low rank and/or sparse matrix for both matrix completion (Jain et al., 2013) and RPCA (Netrapalli et al., 2014; Zhang et al., 2018; Yi et al., 2016). Additionally, both of these problems have been shown to have no spurious local minimizers, and so the ground truth matrices are the only minimizers under some assumptions (Ge et al., 2016) (Ge et al., 2017).

2. Alternating Proximal Gradient Descent

Many different methods have been shown to be effective when minimizing nonconvex relaxations of the l_0 and rank functions, including iterative reweighted methods, and methods based off of low rank factorization. In this paper, we focus on the most commonly used technique: alternating proximal gradient descent.

The proximal gradient descent method applied to the function $F(x) = \phi(x) + g(x)$ iteratively solve the following problem:

$$\begin{aligned} x^{k+1} &= \underset{x}{\operatorname{argmin}} \phi(x) + \frac{1}{2\tau} \|x - (x^k - \tau \nabla g(x^k))\|^2 \\ &:= \text{prox}_{\phi}^\tau(x^k - \tau \nabla g(x^k)) \end{aligned}$$

where we defined the *proximal operator* of a function as the minimum of a combination of the function and the distance from a given point. For many functions ϕ that we are interested in, the proximal operator has a closed form

solution, which we show in Table 1.

For each of the sparsity promoting regularizers in Table 1, the proximal operator is also dubbed as a *shrinkage operator* or a *thresholding operator* because when the input is less than τ , the output is 0. Otherwise, the input is moved towards zero or, for some nonconvex regularizers, is unchanged. So, we can view the proximal gradient algorithm as iteratively taking a step in the gradient direction of $g(x)$, and then applying the proximal operator to promote sparsity.

When applied to the optimization problem in Equation (1), we have

$$\tilde{L}^{k+1} = L^k - \tau_L \frac{d_1 d_2}{n} \mathcal{A}_L^* (\mathcal{A}_L(L^k) + A_s s^k - b) \quad (2a)$$

$$L^{k+1} = \text{prox}_{\phi_{\gamma_L}^{\tau_L \lambda_L}}(\tilde{L}^{k+1}) \quad (2b)$$

$$\tilde{s}^{k+1} = s^k - \tau_S \frac{d_s}{n} A_s^T (\mathcal{A}_L(L^{k+1}) + A_s s^k - b) \quad (2c)$$

$$s^{k+1} = \text{prox}_{\phi_{\gamma_s}^{\tau_s \lambda_s}}(\tilde{s}^{k+1}) \quad (2d)$$

To further simplify the problem, the following proposition will allow L subproblem to be solved in each singular value separately.

Proposition 1. Consider the optimization problem

$$\min_X \sum_i \phi_\gamma(\sigma_i(X)) + \frac{1}{2\tau} \|X - Y\|_F^2 \quad (3)$$

where ϕ_γ is a ν weakly convex function. If $\nu < \tau$, then equation (3) is strongly convex and the minimizer X^* has the same singular vectors as Y with singular values given by

$$\begin{aligned} \sigma_i(X^*) &= \underset{x}{\operatorname{argmin}} \phi_\gamma(x) + \frac{1}{2\tau} (x - \sigma_i(Y))^2 \\ &:= \text{prox}_{\phi_\gamma}^\tau(\sigma_i(Y)) \end{aligned}$$

where $\text{prox}_{\phi_\gamma}^\tau(\sigma_i(Y))$ is the proximal operator.

Proposition 1 tells us that L^k has the singular vectors of \tilde{L}^k and singular values given by

$$\sigma_i(L^k) = \text{prox}_{\phi_\gamma}^{\tau_L \lambda_L}(\sigma_i(\tilde{L}^k)).$$

Likewise, the subproblem in S can be solved in each entry of s individually.

The slowest operation in the alternating proximal gradient method is by far the singular value decomposition. However, in practice we can reduce the number of operations by calculating the truncated singular value decomposition only using the first r_0 singular values, where r_0 is an upper bound on the rank, and enforce that the remaining singular

values are zero. Alternatively, we can calculate each singular value in descending order and stop when a singular value falls below λ_L , as all remaining singular values will be set to zero by the proximal operator. So, in the case of RPCA where each entry is observed, each iteration has a computational complexity of $\mathcal{O}(d_1 d_2 r_0)$, which matches other state of the art methods.

Algorithm 1 Alternating Proximal Gradient Descent for Low-Rank Plus Sparse Optimization

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for  $k = 1, \dots$  do
     $g^{k+1} = -\tau_L \frac{d_1 d_2}{n} \mathcal{A}_L^* (\mathcal{A}_L(U \Sigma^k V^k) + A_s s^k - b)$ 
     $[U^{k+1}, \tilde{\Sigma}^{k+1}, V^{k+1}] = \text{LRSSVD}(U^k, \Sigma^k, V^k, g^{k+1})$ 
     $\Sigma^{k+1} = \text{prox}(\tilde{\Sigma}^{k+1})$ 
     $\tilde{s}^k = s^k - \tau_s \frac{d_s}{n} A_s^T (\mathcal{A}_L(U \Sigma V) + A_s s^k - b)$ 
     $s^{k+1} = \text{prox}^{\lambda_s}(\tilde{s}^{k+1})$ 
end for
    
```

In the case of matrix completion, however, only a sparse set of entries of the low rank matrix are observed, which could be used to increase the efficiency by reducing the amount of computation needed to find the singular value decomposition of L at each iterations. For a low rank matrix with a low rank factorization $U \Sigma V$, we refer to the problem of finding the SVD of the rank r approximation to the matrix $U \Sigma V + g$ for a sparse matrix g as Low Rank plus Sparse SVD (LRSSVD), originally proposed by (Jain et al., 2010).

The LRSSVD task can be accomplished efficiently using the same methods as if we were to find the SVD of any other matrix, such as the Power Iteration method. Recall that the computational complexity of the Power Iteration is limited by the amount of operations needed to multiply the matrix by a vector. Because the computational complexities of calculating both $u(U \Sigma V^T + Y)$ for $u \in \mathbb{R}^{d_1}$ and $(U \Sigma V^T + Y)v$ for $v \in \mathbb{R}^{d_2}$ are $\mathcal{O}((d_1 + d_2)r_0 + n)$, we can calculate the top r_0 singular values and vectors of $X + Y$ with only $\mathcal{O}((d_1 + d_2)r_0^2 + nr_0)$ operations.

The other operations in Algorithm 1 take no more time than the LRPSSVD. The gradient in the L direction, g^k , requires calculating $\Sigma_{ii}^k U_i^k (V_i^k)^T$ for each entry in the support of \mathcal{A}_L . In the case of matrix completion, is nr operations, which matches the computational complexity per iteration for state of the art matrix completion algorithms.

3. Analysis of Proximal Gradient Descent Algorithm

In this section, we present the main result of the paper: a recursive bound on the difference of the iterates of the alternating proximal gradient algorithm and the ground truth low rank matrix and sparse vector. We present the bound for the most general case, and give results on specific prob-

lems in the following section.

3.1. Restricted Isometry and Orthogonality Properties

In order to bound the error in the output of our algorithm relative to the underlying ground truth low-rank and sparse matrices L^* and s^* , we must first make a number of assumptions about L^* , s^* , and the observation models \mathcal{A}_L and A_s .

First, we must assure that the low rank matrix L^* can be separated from a sparse matrix – that is, L^* is not sparse itself. Not only is this necessary for low-rank plus sparse decomposition, but for the problem of matrix completion, this assumption is necessary to assure that a sparse set of observations is a good representation of the entire matrix. For example, consider the matrix consisting of zeros in every entry besides one entry, where the value is 1. We must observe every entry in the matrix to assure that we can reconstruct the matrix exactly, due to the fact that we must observe the nonzero entry and every entry in its row and column.

To exclude such ill-posed problems from our analysis, we will assume that L^* is *incoherent*, as defined as follows.

Definition 2. For a rank r matrix L with singular value decomposition $U \Sigma V^T$, for orthonormal matrices $U \in \mathbb{R}^{d_1 \times r}$, $V \in \mathbb{R}^{d_2 \times r}$, and diagonal matrix $\Sigma \in \mathbb{R}^{r \times r}$, the tangent space of L is defined as

$$\mathcal{T} = \left\{ UX + YV \mid X \in \mathbb{R}^{r \times d_2}, Y \in \mathbb{R}^{d_1 \times r} \right\} \quad (4)$$

Furthermore, we say this tangent space is μ -incoherent if

$$\|U_{:i}\|_2^2 \leq \frac{\mu r}{d_1}, \|V_{:j}\|_2^2 \leq \frac{\mu r}{d_2} \quad \forall i, j \quad (5)$$

We define the projection of a matrix onto the sparse space Ω as

$$\mathcal{P}_\Omega(X) = \begin{cases} X_{ij} & (i, j) \in \Omega \\ 0 & \text{else} \end{cases} \quad (6)$$

and onto the tangent space \mathcal{T} as

$$\mathcal{P}_\mathcal{T}(X) = UU^T X + X V V^T - UU^T X V V^T \quad (7)$$

Next, we we discuss the conditions that the observation models \mathcal{A}_L and A_s must satisfy in order to recover the ground truth low rank and sparse matrix, known as the restricted isometry property. Loosely, the RIP states that for any two vectors in Ω (or matrices \mathcal{T}), we can obtain a sufficiently accurate estimate of the distance between the two through the observation model A_s (or \mathcal{A}_L).

Definition 3. The linear mapping A_s satisfies the (α, κ) sparse Restricted Isometry Property if, for any x satisfying

$$\|x\|_0 \leq \alpha d_s$$

$$(1 - \kappa_S)\|x\|^2 \leq \tau_s \frac{n}{d_s} \|A_s x\|^2 \leq (1 + \kappa_S)\|A_s x\|^2$$

for some constant τ_s . Likewise, the linear mapping \mathcal{A} satisfies the (μ, r, κ) low rank Restricted Isometry Property if for any X in a μ -incoherent rank r tangent space \mathcal{T} ,

$$(1 - \kappa_L)\|X\|_F^2 \leq \tau_L \frac{n}{d_1 d_2} \|\mathcal{A}_L X\|_F^2 \leq (1 + \kappa_L)\|X\|_F^2$$

for some constant τ_L .

In some cases, it may be more useful to use the following characterization of the RIP, which bounds the difference between the operator $\tau \mathcal{A}^* \mathcal{A}$ and the identity operator when restricted to sparse vectors or low-rank and incoherent matrices.

Proposition 2. For a matrix $A_s \in \mathbb{R}^{n \times d_s}$ satisfying the (α, κ_s) sparse RIP,

$$\left\| \frac{\tau_s n}{d_s} \mathcal{P}_\Omega A_s^T A_s \mathcal{P}_\Omega - \mathcal{P}_\Omega \right\|^2 \leq \kappa_s.$$

Likewise, for any linear mapping $\mathcal{A}_L : \mathbb{R}^{d_1 \times d_2} \rightarrow \mathbb{R}^n$ satisfying the (μ, r, κ_L) low rank RIP,

$$\left\| \frac{\tau_L n}{d_1 d_2} \mathcal{P}_\mathcal{T} \mathcal{A}_L^* \mathcal{A}_L \mathcal{P}_\mathcal{T} - \mathcal{P}_\mathcal{T} \right\|^2 \leq \kappa_L$$

Finally, we discuss the interplay between the set of sparse matrices and the low rank, incoherent tangent space, and their observation models. We hope to be able to separate the measurement vector b into two parts: one in the span of $\mathcal{A}_L \mathcal{P}_\mathcal{T}$, and one in the span of $A_s \mathcal{P}_\Omega$. In order for this to be able to be done quickly, we require there to be no non-trivial vectors in the intersection of the two sets, which is equivalent to saying that $\|\mathcal{P}_\Omega A_s^T \mathcal{A}_L \mathcal{P}_\mathcal{T}\| < 1$. Under some assumptions, this norm is actually close to zero, a concept we refer to as restricted orthogonality, which we define here and verify applies to the problems we are interested in in Section 4.

Definition 4. Linear maps \mathcal{A}_L and A_s satisfy the restricted orthogonality property over the sets \mathcal{T} and Ω (respectively) when

$$\|\mathcal{P}_\Omega A_s^* \mathcal{A}_L \mathcal{P}_\mathcal{T}\|^2 \leq \kappa_{SL}$$

$$\|\mathcal{P}_\mathcal{T} \mathcal{A}_L^* A_s \mathcal{P}_\Omega\|^2 \leq \kappa_{LS}$$

3.2. Main Result

Define the difference between the iterates of the alternating proximal gradient descent algorithm and the ground truth low rank matrix and sparse vector at iteration k as $\Delta_L^k = L^* - L^k$ and $\Delta_s^k = s^* - s^k$. Our main result in the most general form gives a bound on the norm of Δ_L^k and Δ_s^k in terms of the differences at the previous iteration, Δ_L^{k-1} and Δ_s^{k-1} .

Theorem 1. Let L^k and s^k be the sequences generated by Algorithm 1. Assume that

$$b = \mathcal{A}_L(L^*) + A_s s^* + \mathcal{E} \in \mathbb{R}^n,$$

where $L^* \in \mathbb{R}^{d_1 \times d_2}$ is a rank r and μ -incoherent matrix, and $s^* \in \mathbb{R}^{d_s}$ is a sparse vector with $\text{supp}(s^*) = \Omega$, and the linear mappings \mathcal{A}_L and A_s satisfy the $(2r, 3\mu, \kappa_L)$ -low rank RIP and the (α, κ_s) -sparse RIP respectively, and together satisfy the ROP with constant κ_{LS} . If $\lambda_L \geq \|\mathcal{A}_L^*(\mathcal{E})\|_2 + \|\mathcal{A}_L^* A_s \Delta_s^{k-1}\|_2$, then

$$\begin{aligned} \|\Delta_L^{k+1}\|_F^2 &\leq \kappa_L \|\Delta_L^k\|_F^2 + \kappa_{LS} \frac{\tau_L d_1 d_2}{n} \|\Delta_s^k\|_F^2 \\ &\quad + \frac{\tau_L d_1 d_2}{n} \|\mathcal{P}_\mathcal{T} \mathcal{A}_L^* \mathcal{E}\|_F^2 + \lambda_L r \phi'(\sigma_r(L^*) - 2\lambda_L) \end{aligned}$$

Likewise, if $\lambda_S \geq \|A_s^T(\mathcal{E})\|_\infty + \|A_s^T \mathcal{A}_L \Delta_L^{k-1}\|_\infty$, then

$$\begin{aligned} \|\Delta_s^{k+1}\|_F^2 &\leq \kappa_S \|\Delta_s^k\|_F^2 + \frac{\tau_s d_s}{n} \kappa_{LS} \|\Delta_L^{k+1}\|_F^2 \\ &\quad + \frac{\tau_s d_s}{n} \|\mathcal{P}_\Omega A_s^T \mathcal{E}\|_F^2 + \lambda_s \alpha d_s \phi'(s_{\min} - 2\lambda_s) \end{aligned}$$

For each of these bound, we can think of the first term as the estimation error introduced by the noise, and the second term as the approximation error, which accounts for the bias in the regularizer proportional to the derivative of the regularizer. Previous result for the nuclear norm and l_1 norm give similar bounds, but make the concession that the approximation error is the dominating term. Under some circumstances, that term is equal to zero in our bound.

3.3. Proof of Main Result

We start by presenting the the following two lemmas regarding the proximal operator for the sparse and low rank regularizers.

Lemma 1. Let L^* be a rank r , μ -incoherence matrix whose singular vectors form the tangent space \mathcal{T} , and \bar{L} be defined as

$$\bar{L} = \underset{L \in \mathbb{R}^{d_1 \times d_2}}{\text{argmin}} \lambda \Phi(L) + \|L - L^* + \delta\|_F^2 \quad (8)$$

Where Φ is an at most $\frac{1}{\lambda}$ weakly convex regularizer satisfying Assumption 1, and $\delta \in \mathbb{R}^{d_1 \times d_2}$ satisfies $\|\delta\|_2 \leq \lambda$. Define $\Delta_L = \bar{L} - L^*$. Then,

$$\|\Delta_L\|_F^2 \leq \|\mathcal{P}_\mathcal{T}(\delta)\|_F^2 + 2\lambda \sqrt{r} \phi'(\sigma_r(L^*)) \quad (9)$$

A proofs of Lemma 1 can be found in Appendix A.

In order to utilize the RIP and ROP conditions, we need to verify that L^{k+1} is low rank and incoherent, and that s^{k+1} is sparse, which we will do inductively. Assume that L^k is at most rank r , and that its tangent space is 2μ incoherent. Additionally, assume that $\text{supp}(S^k) \subseteq \text{supp}(S^*)$. Clearly,

these conditions are met at the first iteration by initializing the algorithm with $L^0 = 0$ and $S^0 = 0$.

At iteration $k + 1$,

$$\begin{aligned} L^* - \bar{L}^{k+1} &= L^* - \left(L^k - \frac{\tau_L n}{d_1 d_2} \mathcal{A}_L^* (\mathcal{A}_L(L^k) + A_s(S^k) - b) \right) \\ &= L^* - L^k + \frac{\tau_L n}{d_1 d_2} \mathcal{A}_L^* \mathcal{A}_L L^k + \frac{\tau_L n}{d_1 d_2} \mathcal{A}_L^* A_s S^k \\ &\quad - \frac{\tau_L n}{d_1 d_2} \mathcal{A}_L^* (\mathcal{A}_L(L^*) + A_s(S^*) + \mathcal{E}) \\ &= (\Delta_L^k - \frac{\tau_L n}{d_1 d_2} \mathcal{A}_L^* \mathcal{A}_L \Delta_L^k) - \frac{\tau_L n}{d_1 d_2} (\mathcal{A}_L^* A_s \Delta_S^k - \mathcal{A}_L^* \mathcal{E}) \end{aligned}$$

where the first equality comes from the definition of b , and the second inequality substitutes $\Delta_S^k = S^* - S^k$ and $\Delta_L^k = L^* - L^k$. By Lemma 1 (along with the triangle inequality), we have that

$$\begin{aligned} \|\Delta_L^{k+1}\|_F^2 &\leq \|\mathcal{P}_T(\Delta_L^k - \tau_L \frac{d_1 d_2}{n} \mathcal{A}_L^* \mathcal{A}_L \Delta_L^k)\|_F^2 \\ &\quad + \tau_L \frac{d_1 d_2}{n} \|\mathcal{P}_T \mathcal{A}_L^* A_s \Delta_S^k\|_F^2 + \tau_L \frac{d_1 d_2}{n} \|\mathcal{P}_T \mathcal{A}_L^* \mathcal{E}\|_F^2 \\ &\quad + \lambda_L r \phi'(\sigma_r(L^*) - 2\lambda_L \tau_L) \end{aligned}$$

Let \mathcal{T}^k denote the union of the tangent space of the rank r approximation L^k and \mathcal{T} so that $\Delta_L^k \in \mathcal{T}^k$. Then,

$$\begin{aligned} \|\mathcal{P}_T(\Delta_L^k - \tau_L \frac{d_1 d_2}{n} \mathcal{A}_L^* \mathcal{A}_L \Delta_L^k)\|_F^2 &\leq \|\mathcal{P}_{\mathcal{T}^k}(\Delta_L^k - \tau_L \frac{d_1 d_2}{n} \mathcal{A}_L^* \mathcal{A}_L \Delta_L^k)\|_F^2 \\ &\leq \|\mathcal{P}_{\mathcal{T}^k}(\tau_L \frac{d_1 d_2}{n} \mathcal{A}_L^* \mathcal{A}_L - \mathcal{I}) \mathcal{P}_{\mathcal{T}^k} \Delta_L^k\|_F^2 \leq \kappa_L \|\Delta_L^k\|_F^2 \end{aligned}$$

where the first inequality comes from the contractive property of $\mathcal{P}_{\mathcal{T}}$, and the second inequality comes from the fact that $\mathcal{P}_{\mathcal{T}^k} \Delta_L^k = \Delta_L^k$. And, because \mathcal{T}^k is has incoherence at most 3μ , we can apply the low-rank RIP to obtain the third inequality.

By the inductive hypothesis stating that $\text{supp}(S^k) \subseteq \text{supp}(S^*)$ (and thus, that S^k is α -sparse), we can use the ROP to claim that

$$\|\mathcal{P}_T \mathcal{A}_L^* A_s \Delta_S^k\|_F^2 \leq \kappa_{SL} \|\Delta_S^k\|_F^2.$$

Combining these gives the desired bound on Δ_L^k .

Now, we must show that L^{k+1} is rank r and 2μ incoherent. By the tk inequality (stated in Appendix C), we know that

$$\begin{aligned} \sigma_{r+1}(\bar{L}^{k+1}) &\leq \sigma_{r+1}(L^*) + \|L^* - \bar{L}^{k+1}\|_2 \\ &\leq \|\Delta_S^k\|_2 + \|\mathcal{A}_L^* \mathcal{E}\|_2 \end{aligned}$$

Because this is less than λ_L by our assumptions, L^{k+1} must be rank r .

In order to show that \bar{L}^{k+1} is at most 2μ incoherent, we apply the following Lemma, which utilizes the David-Kahan inequality.

Lemma 2. *If X is a rank r , μ -incoherent matrix, and Δ satisfies $\|\Delta\|_2 \leq \frac{1}{4}\sigma_r(X)$, then the top r singular vectors of the matrix $X + \Delta$ form a 2μ -incoherent tangent space.*

A proof of Lemma 2 can be found in Appendix A. We conclude the proof by bounding Δ_s^k and showing $s^{k+1} \in \Omega$ in a similar manner in Appendix A.

4. Results for Specific Models

In this section, we utilize Theorem 1 to analyze an application of the alternating proximal gradient descent algorithm to the problems of matrix completion and RPCA.

4.1. Matrix Completion

We start by considering the problem of matrix completion. Here, we have a sparse set of observed entries of L , $\mathcal{A}_L = \mathcal{A}_\Omega$, and we do not consider a sparse vector s (i.e. $A_s = 0$).

We present a version of the RIP for the sampling operator from (Recht, 2011).

Lemma 3 ((Recht, 2011)). *Let Ω be a set of n entries of $\{1, \dots, d_1\} \times \{1, \dots, d_2\}$ drawn independently at random with uniform probability, with $n > 64\mu r(d_1 + d_2) \log(d_2)$. Then, with probability at least $1 - 2d_2^{-2}$,*

$$\frac{5}{6} \|X\|_F^2 \leq \frac{d_1 d_2}{n} \|\mathcal{A}_\Omega(X)\|^2 \leq \frac{7}{6} \|X\|_F^2$$

for any rank r , μ -incoherent matrix X .

We now present an error bound of for a stationary point of Algorithm 1 applied to matrix completion.

Theorem 2. *Let $b = \mathcal{A}_\Omega X^* + \mathcal{E}$ for a rank r , μ -incoherent matrix $X^* \in \mathbb{R}^{d_1 \times d_2}$. Under the same assumptions as Lemma 3, if $\lambda > \|\mathcal{A}_\Omega^*(\mathcal{E})\|_2$, then the iterates of Algorithm 1 linearly converge to a point \bar{X} such that $\|X^* - \bar{X}\|_F^2$ is less than:*

$$\frac{6}{5} \frac{d_1 d_2}{n} \left(\underbrace{\|\mathcal{P}_T \mathcal{A}_\Omega^* \mathcal{E}\|_F^2}_{\text{optimal error rate}} + \underbrace{\lambda r \phi'(\sigma_{r+1}(X^*) - 2\lambda)}_{\text{bias term}} \right)$$

with convergence rate $\frac{1}{6} \frac{d_1 d_2}{n}$.

A proof of Theorem 2 can be found in Appendix B. The two terms of the error bound account for the optimal error rate and a bias term. The optimal error rate is the error bound if we know the tangent space of X^* a priori, that is, the difference between X^* and the solution to the optimization problem

$$\min_{X \in \mathcal{T}} \|\mathcal{A}_\Omega(X) - b\|_F^2.$$

The oracle rate is further discussed in (Candes & Plan, 2010) and (Negahban & Wainwright, 2012).

4.2. Robust PCA

Next, we will use Theorem 1 to analyze the APGD algorithm applied to the problem of RPCA. Specifically, we are interested in the special case of (1) where the $A_s = I_n$, and $A_L = A_{\Omega^{obs}}$.

In order for RPCA to be possible, we need the nonzero entries of $A_{\Omega^{obs}}^* s^*$ to be sufficiently well-distributed throughout the rows and columns – if the sparse corruptions affected the same row or column of L^* , then this would also be a low-rank perturbation and thus be impossible to separate from L^* without further information. So, we will assume that $A_L^* s^*$ is α -sparse, defined as follows.

Definition 5. *The matrix S is α -sparse for $0 < \alpha < 1$ if the proportion of nonzero entries in any row or column is less than α . That is,*

$$\|S_{i:}\|_0 \leq \alpha d_1, \|S_{:j}\|_0 \leq \alpha d_2 \quad \forall i, j \quad (10)$$

In order to verify that the ROP property holds, we present the following Lemma:

Lemma 4. *Let \mathcal{T} be a rank r , μ -incoherent tangent space, and let Ω be an α -sparse subspace. Then,*

$$\|\mathcal{P}_{\mathcal{T}} \mathcal{P}_{\Omega}\| \leq 2\alpha\mu r, \|\mathcal{P}_{\Omega} \mathcal{P}_{\mathcal{T}}\| \leq 2\alpha\mu r \quad (11)$$

We can now give a bound for the stationary point of the APGD algorithm applied to RPCA.

Theorem 3. *Let $b = A_{\Omega^{obs}}(L^*) + s^* + \mathcal{E}$ for a rank r , μ -incoherent matrix $L^* \in \mathbb{R}^{d_1 \times d_2}$, and a sparse vector $s^* \in \mathbb{R}^n$ with $\|s^*\|_{\infty} \leq 2\|L^*\|_{\infty}$. Under the same assumptions on Ω^{obs} as Lemma 3, and assuming that $A_{\Omega}^* s^*$ is α -sparse and $\alpha\mu r \leq \frac{1}{64}$. If $\lambda_L \geq \frac{1}{6} + \|A_{\Omega}^* \mathcal{E}\|_2$ and $\lambda_s \geq \frac{\mu r}{d_1} + \|\mathcal{E}\|_{\infty}$, then the iterates of Algorithm 1 linearly converge to a point \bar{L}, \bar{s} satisfying*

$$\begin{bmatrix} \|L^* - \bar{L}\|_F^2 \\ \|s^* - \bar{s}\|_2^2 \end{bmatrix} \leq \frac{6}{5} \begin{bmatrix} 1 \\ \frac{1}{64} \end{bmatrix} \begin{bmatrix} \frac{d_1 d_2}{64^n} \\ \frac{5}{6} \end{bmatrix} \begin{bmatrix} \frac{d_1 d_2}{n} \|\mathcal{P}_{\mathcal{T}} A_L^* \mathcal{E}\|_F^2 \\ \|\mathcal{P}_{\Omega} A_s^T \mathcal{E}\|_F^2 \end{bmatrix}$$

with convergence rate $\frac{1}{6} \frac{d_1 d_2}{n}$.

5. Numerical Results

We implemented Algorithm 1 in both Matlab R2020a, for which the code is available at GitHub. We also made the code available as part of the Python package SpaLoR. All results in this section are obtained with the Matlab version in order to accurately compare to other algorithms which are only available in Matlab, and are run on a Windows 10 desktop with an AMD Phenom 3.40 GHz processor and 8 Gb of RAM.

5.1. Matrix Completion

We compare our method for matrix completion to another method utilizing nonconvex regularizer from (Yao et al., 2019) (FaNCL), along with a method to minimize the nuclear norm IALM, from (Lin et al., 2010), and a rank constrained method, LMaFit (Wen et al., 2012). The results are shown in Table 2.

We start by comparing the performance of the methods on randomly generated low rank matrices of varying size, rank, percentage of observed entries, and standard deviation of the noise in the measurements (shown relative to the mean absolute value of the low rank matrix). In each of the cases, our method performs exactly as well as FaNCL and LMaFit when the correct rank is given. IALM performs equally well in the first case, and slightly worse than the remaining three cases due to the fact that the nuclear norm biases the result towards zero.

Next, we show the results on common data sets for recommendation systems, the Jester data set (Goldberg et al., 2001) and MovieLens 1M (mov). For each of these two data sets, we partition the observations into five folds, fit a low rank model to four of the folds and calculate the accuracy on the remaining fold. We do this for each of the five folds and present the average normalized mean absolute error. In both cases, our method outperforms the other three algorithms we compare to.

5.2. Robust PCA

We compare our method to several other prominent RPCA methods, including LMaFit (Shen et al., 2014), AltProj (Netrapalli et al., 2014), RPCA-GD (Yi et al., 2016), and IALM (Lin et al., 2010). We compared with many other methods included in the LRSLibrary (Sobral et al.), however we only include results from the aforementioned algorithms as they gave the most accurate results for matrices a significant amount with Gaussian noise, a test case we emphasise in this section.

It is worth noting that, while our algorithm does not require an estimate of the rank of L^* or the sparsity of S^* a priori, BLANK requires both, and LMaFit and AltProj require an estimate of the rank. However, we found that LMaFit and AltProj still perform very well when this estimate is unreliable, as LMaFit includes a rank-estimation scheme and AltProj starts by performing a rank 1 projection, and increases the rank up until the estimate given. We provide all methods with an upper bound on the rank equal to twice the rank of L^* and an upper bound on the number of corrupted entries equal to twice that of S^* .

The most commonly utilized test case in the literature for RPCA is the task of separating the background and foreground of a video. In this scenario, each frame of the video

Table 2. Comparison of four different matrix completion algorithms on randomly generated low rank matrices and common recommendation data sets. The algorithm LMaFit reconstructs a matrix of a given rank k . The table shows the results when the algorithm is given the exact rank ($k = r$) and an incorrect rank ($k = 2r$). Time is given in seconds

Jester Dataset									ML 1M			
d_1	1000		1000		5000		5000		24983		6040	
d_2	500		500		1000		1000		100		3952	
r	5		5		10		10		-			
n/d_1d_2	0.3		0.1		0.2		0.05		0.58		0.034	
$\text{std}(\mathcal{E})$	0.1		0.02		0.1		0.02		-			
	RFNE	T	RFNE	T	RFNE	T	RFNE	T	NMAE	T	NMAE	T
APGD	3.28e-4	0.7	2.90e-4	1.1	1.69e-4	10	1.96e-4	29	0.159	21	0.172	172
FaNCL	3.28e-4	1.8	2.90e-4	4.3	1.69e-4	31	2.49E-4	57	0.183	42	0.200	42
IALM	3.28e-4	2.6	2.92e-4	2.8	1.72e-4	32	1.99e-4	27	0.163	77	0.183	216
LMaFit (2r)	4.68e2	21	1.20e3	6.2	4.99e2	197	1.60e3	40	-			
LMaFit(r)	3.28e-4	0.4	2.90e-4	0.3	1.69e-4	3.9	1.96e-4	3.9	0.168	7.4	9.174	44

Table 3. Accuracy and run-time for RPCA on four videos of fish.

	Marine Snow		Small Aquaculture		Caustics		Two Fish		Fish Swarm	
	Acc	Time (s)	Acc	Time (s)	Acc	Time (s)	Acc	Time (s)	Acc	Time (s)
APGD	0.92	87	0.92	592	0.88	459	0.90	467	0.75	473
LMaFit	0.59	258	0.65	316	0.7	273	0.62	291	0.55	310
AltProj	0.84	143	0.62	109	0.86	83	0.76	109	0.62	130
RPCA-GD	0.92	94	0.72	77	0.87	76	0.96	96	0.66	93
IALM	0.57	379	0.56	286	0.56	288	0.56	342	0.56	305

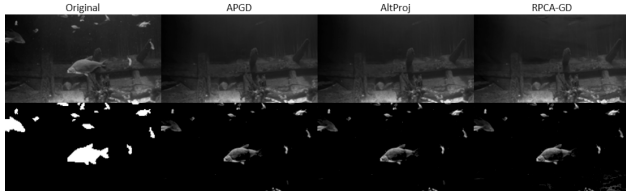


Figure 1. Frames from the Marine Snow video. The first row of the first column shows the original frame, with the image segmentation for that frame below it. The remaining three columns show the background and foreground obtained by three different methods.

is represented as a column vector in M . If the background is static (or, at least, in some way repetitive), then we can expect the background of each frame to be represented as a low rank matrix, and the foreground as a sparse matrix. See (Candès et al., 2011) for further details.

We evaluated our algorithm on the Underwater Change Detection dataset (Radolko et al., 2016). Out of the 1100 frames in each video, the ground truth image segmentation is included for last 100 frames. This allows use to present an objective and accurate metric of how well each method is able to identify the foreground of the image.

In Table 5.2, we present the runtime and the accuracy of

determining which pixels contain a fish for our method compared to the four previously mentioned approaches. To calculate the accuracy, we average of the true positive ratio and true negative ratio. In three of the five videos, our method achieves the highest accuracy, where as in the other two the RPCA-GD algorithm preforms slightly better. The recovered background and foreground for our method, RPCA-GD and AltProj are shown in Figure 1, along with the original frame and segmented image. Additional images are shown in Appendix D.

6. Conclusions

We have shown an novel convergence analysis of the alternating proximal gradient descent algorithm applied to the problems of matrix completion and RPCA with nonconvex regularizers, and bound the difference from the ground truth low rank matrix and sparse vector. Future work on the topic could include extending our analysis to data that lies on more complicated, nonlinear manifolds.

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