

A General Framework for Sparsity Regularized Feature Selection via Iteratively Reweighted Least Square Minimization

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Abstract

A variety of feature selection methods based on sparsity regularization have been developed with different loss functions and sparse regularization functions. Capitalizing on the existing sparsity regularized feature selection methods, we propose a general sparsity feature selection (GSR-FS) algorithm that optimizes a $\ell_{2,r}$ -norm ($0 < r \leq 2$) based loss function with a $\ell_{2,p}$ -norm ($0 < p \leq 1$) sparse regularization. The $\ell_{2,r}$ -norm ($0 < r \leq 2$) based loss function brings flexibility to balance data-fitting and robustness to outliers by tuning its parameter, and the $\ell_{2,p}$ -norm ($0 < p \leq 1$) based regularization function is able to boost the sparsity for feature selection. To solve the optimization problem with multiple non-smooth and non-convex functions when $r, p < 1$, we develop an efficient solver under the general umbrella of Iterative Reweighted Least Square (IRLS) algorithms. Our algorithm has been proved to converge with a theoretical convergence order of at least $\min(2 - r, 2 - p)$. The experimental results have demonstrated that our method could achieve competitive feature selection performance on publicly available datasets compared with state-of-the-art feature selection methods, with reduced computational cost.

1. Introduction

Feature selection plays an important role in high-dimensional data analysis for selecting informative features and removing irrelevant or redundant ones (Cawley et al., 2006; Guyon & Elisseeff, 2003; Kira & Rendell, 1992; Lewis, 1992; Peng et al., 2005). Among existing feature selection methods, sparsity regularization based methods are appealing for their excellent performance (Argyriou & Evgeniou, 2007; Bradley & Mangasarian, 1998; Liu et al., 2009; Nie et al., 2010; Obozinski et al., 2006; Tibshirani, 1996; Wang et al., 2008; Xiang et al., 2012). In particular, ℓ_1 -norm has been widely adopted in feature selection algorithms, such as

Lasso (Tibshirani, 1996) and sparse SVM (Bradley & Mangasarian, 1998; Wang, et al., 2008). Built upon ℓ_1 -norm based regularization models, $\ell_{2,1}$ -norm has been used for feature selection in problems with multiple tasks or multiple classes (Argyriou & Evgeniou, 2007; Liu, et al., 2009; Nie, et al., 2010; Obozinski, et al., 2006; Xiang, et al., 2012). More recently, ℓ_p -norm and $\ell_{2,p}$ -norm ($0 < p < 1$) based regularization models have gained increasing attention (Bolon-Canedo, et al., 2013; Chartrand & Staneva, 2008; Kong & Ding, 2014; Liu et al., 2007; Peng & Fan, 2016; Zhang et al., 2014) since they can yield sparser solutions than ℓ_1 -norm and $\ell_{2,1}$ -norm based models (Chartrand, 2007; Zeng et al., 2014).

Although a variety of sparsity regularization based feature selection methods with different sparse regularization functions have been developed, most of them adopt a least square loss function. The least square loss function has good data-fitting performance. However, it is sensitive to outliers. A robust feature selection (RFS) method with joint $\ell_{2,1}$ -norm minimization on both the loss function and regularization function was proposed (Nie, et al., 2010; Xiang, et al., 2012) and has been extended (Wang & Chen, 2013) with joint $\ell_{2,p}$ -norm ($0 < p \leq 1$). However, it is not necessary to use the same norm for both the loss function and sparse regularization function. To make the sparsity regularized feature selection method more flexible, we propose a general sparsity regularized feature selection (GSR-FS) algorithm that optimizes a $\ell_{2,r}$ -norm ($0 < r \leq 2$) based loss function and a $\ell_{2,p}$ -norm ($0 < p \leq 1$) sparse regularization function. Particularly, the $\ell_{2,r}$ -norm ($0 < r \leq 2$) based loss function can balance the data fitting and robustness to outliers and the $\ell_{2,p}$ -norm ($0 < p < 1$) based regularization function is able to boost the model sparsity for feature selection.

The optimization algorithms used in the existing sparsity regularized methods typically handle optimization problems

with one non-smooth term¹ and are not suitable for our optimization problem with two non-smooth terms when $r, p \leq 1$. Iteratively reweighted least squares (IRLS) based methods have been widely used to solve sparse optimization problems in many fields (Candes et al., 2008; Chartrand & Yin, 2008; Gorodnitsky & Rao, 1997; Lu et al., 2014). However, the existing IRLS based algorithms only handle optimization problems with no more than one non-smooth function (Lu et al., 2015). To optimize our problem, we develop a novel algorithm based on IRLS with a convergence order of at least $\min(2 - r, 2 - p)$.

Our method has been validated based on 6 publicly available datasets and achieved competitive feature selection performance with respect to both classification accuracy and computational cost compared with 6 state-of-the-art feature selection algorithms, including Minimum-Redundancy Maximum-Relevance (mRMR) (Peng, et al., 2005), ReliefF (Kira & Rendell, 1992), Multi-Task Feature Selection (MTFS) (Argyriou & Evgeniou, 2007; Liu, et al., 2009; Obozinski, et al., 2006), Robust Feature Selection (RFS) (Nie, et al., 2010; Xiang, et al., 2012), an extended RFS (E-RFS) (Wang & Chen, 2013), and Rank One Update Algorithm (RK1U) (Zhang, et al., 2014).

2.A unified sparse feature selection algorithm

Given a matrix $\mathbf{A} \in \mathbb{R}^{m \times n}$, its $\ell_{2,p}$ -norm ($r > 0, p > 0$) is defined as:

$$\|\mathbf{A}\|_{2,p} = \left(\sum_{i=1}^m \left(\sum_{j=1}^n |\mathbf{a}_{i,j}|^2 \right)^{\frac{p}{2}} \right)^{\frac{1}{p}} = \left(\sum_{i=1}^m (\|\mathbf{a}_i\|_2^2)^{\frac{p}{2}} \right)^{\frac{1}{p}}, \quad (1)$$

where $\|\mathbf{a}_i\|_2$ denotes ℓ_2 -norm of the i -th row vector of \mathbf{A} .

Given m training samples $\mathbf{X} = \{\mathbf{x}^i\}_{i=1}^m$, $\mathbf{x}^i \in \mathbb{R}^n$, belonging to c ($c \geq 2$) classes, and their class labels $\mathbf{Y} = \{\mathbf{f}^i\}_{i=1}^m$, $\mathbf{f}^i = [-1, \dots, 1, \dots, -1] \in \mathbb{R}^c$ (the j -th element is 1 and others are -1 for the i -th data point belonging to the j -th class). In this paper, we adopt a $\ell_{2,r}$ -norm ($0 < r \leq 2$) based loss function and a $\ell_{2,p}$ -norm ($0 < p \leq 1$) based regularization function for feature selection, *i.e.*,

$$\min_{\mathbf{W}} \mathcal{J}(\mathbf{W}) = \|\mathbf{X}\mathbf{W} - \mathbf{Y}\|_{2,r}^r + \lambda \|\mathbf{W}\|_{2,p}^p, \quad (2)$$

where $\mathbf{W} \in \mathbb{R}^{n \times c}$ is the weight matrix to be learned, and non-zero rows of \mathbf{W} indicate the selected features.

We choose $\|\mathbf{X}\mathbf{W} - \mathbf{Y}\|_{2,r}^r$ ($0 < r \leq 2$) instead of the traditional $\|\mathbf{X}\mathbf{W} - \mathbf{Y}\|_F^2$ as the loss function for following reasons. In general, a loss function with smaller r is more robust to outliers, whereas with larger r has better data-fitting performance. As indicated by the plots shown in Figure 1 (a), a small $r \leq 1$ for ℓ_r -norm could reduce the impact of an outlier on the loss function compared with a larger r . The impact of outliers in classification is also illustrated by the simple 2D linear classification models with different settings of ℓ_r -norm based loss functions. In particular, as shown in Figure 1 (b), the classification lines remain unchanged with r if no outlier is present in the training data. However, the classification model with a ℓ_2 -norm based loss function could change dramatically with outliers, and the classification models with a smaller r are relatively more robust to outliers, as illustrated by Figure 1 (c). The regularization function $\|\mathbf{W}\|_{2,p}^p$ has a direct impact on the solution's sparsity, and small $p \leq 1$ is able to boost sparsity.

The proposed method in Eqn. (2) is a generalization of existing sparsity regularization based feature selection methods, and many of them are special cases of the proposed method. Differences between our method and the existing methods under comparison are summarized in Table 1.

3.A novel IRLS method

The optimization problem of Eqn. (2) is a difficult problem with 2 non-convex, non-smooth optimization functions when $0 < r < 1$ and $0 < p < 1$. To efficiently solve this problem, we propose an iterative algorithm, and at each iteration step we remodel the optimization problem of Eqn. (2) as are-weighted least square minimization problem with analytical solutions.

When $\|\mathbf{x}_i\mathbf{W} - \mathbf{y}_i\| \neq 0$ and $\|\mathbf{w}_i\| \neq 0$ ($\mathbf{x}_i\mathbf{W} - \mathbf{y}_i$ and \mathbf{w}_i are the i -th row vector of $\mathbf{X}\mathbf{W} - \mathbf{Y}$ and \mathbf{W} , respectively),

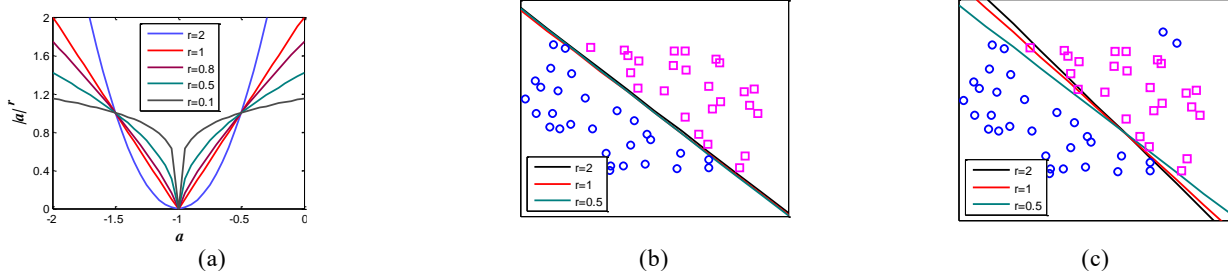


Figure 1: (a) $|a|^r$ with different r . (b) Optimal classification lines for ℓ_r -norm loss functions without any outlier. (c) Optimal classification lines for ℓ_r -norm loss functions with outliers.

¹ RFS and the extended RFS reformulated their optimization objective function with two non-smooth terms as a problem with one non-smooth

term due to both the loss function and the regularization term share the same $\ell_{2,p}$ -norm ($0 < p \leq 1$) form (Nie, et al., 2010; Wang & Chen, 2013).

Table1: Comparisons of MTFs, RK1U, RFS, Extended RFS and our method

Method name	Objective function	Value of r	Value of p	Are r and p must be the same
MTFS	$\ \mathbf{X}\mathbf{W} - \mathbf{Y}\ _F^2 + \lambda\ \mathbf{W}\ _{2,1}$	$r = 2$	$p = 1$	NO
RK1U	$\ \mathbf{X}\mathbf{W} - \mathbf{Y}\ _F^2 + \lambda\ \mathbf{W}\ _{2,p}^p$	$r = 2$	$0 < p \leq 1$	NO
RFS	$\ \mathbf{X}\mathbf{W} - \mathbf{Y}\ _{2,1} + \lambda\ \mathbf{W}\ _{2,1}$	$r = 1$	$p = 1$	YES
E-RFS	$\ \mathbf{X}\mathbf{W} - \mathbf{Y}\ _{2,p}^p + \lambda\ \mathbf{W}\ _{2,p}^p$	$0 < r \leq 1$	$0 < p \leq 1$	YES
Ours	$\ \mathbf{X}\mathbf{W} - \mathbf{Y}\ _{2,r}^r + \lambda\ \mathbf{W}\ _{2,p}^p$	$0 < r \leq 2$	$0 < p \leq 1$	NO

the gradient of $\mathcal{J}(\mathbf{W})$ in Eqn. (2) with respect to \mathbf{W} is

$$\frac{\partial \mathcal{J}(\mathbf{W})}{\partial \mathbf{W}} = 2\mathbf{X}^T \mathbf{S}_1 (\mathbf{X}\mathbf{W} - \mathbf{Y}) + 2\lambda \mathbf{S}_2 \mathbf{W}, \quad (3)$$

where $\mathbf{S}_1 \in \mathbb{R}^{m \times m}$ and $\mathbf{S}_2 \in \mathbb{R}^{n \times n}$ are diagonal matrices, and their i -th diagonal elements are $\mathbf{s}_1^{ii} = r/(2\|\mathbf{x}_i \mathbf{W} - \mathbf{y}_i\|_2^{2-r})$ and $\mathbf{s}_2^{ii} = p/2\|\mathbf{w}_i\|_2^{2-p}$. When $\|\mathbf{x}_i \mathbf{W} - \mathbf{y}_i\| = 0$ and $\|\mathbf{w}_i\| = 0$, we just handle it as the same as that in paper (Nie, et al., 2010).

Setting $\partial \mathcal{J}(\mathbf{W})/\partial \mathbf{W}$ to be 0, then we have a solution of Eqn. (2), i.e.,

$$\mathbf{W} = (\mathbf{X}^T \mathbf{S}_1 \mathbf{X} + \lambda \mathbf{S}_2)^{-1} \mathbf{X}^T \mathbf{S}_1 \mathbf{Y}. \quad (4)$$

If \mathbf{S}_1 and \mathbf{S}_2 are fixed, we can construct an auxiliary objective function $\mathcal{J}_1(\mathbf{W})$ to have the same gradient as $\mathcal{J}(\mathbf{W})$ in Eqn. (2),

$$\mathcal{J}_1(\mathbf{W}) = \|\mathbf{S}_1(\mathbf{X}\mathbf{W} - \mathbf{Y})\|_F^2 + \lambda\|\mathbf{S}_2 \mathbf{W}\|_F^2, \quad (5)$$

where $\mathbf{\Sigma}_1 = (\mathbf{S}_1)^{1/2}$ and $\mathbf{\Sigma}_2 = (\mathbf{S}_2)^{1/2}$, and their i -th diagonal elements are $\mathbf{\zeta}_1^{ii} = \sqrt{r/2}/\|\mathbf{x}_i \mathbf{W} - \mathbf{y}_i\|_2^{1-r/2}$ and $\mathbf{\zeta}_2^{ii} = \sqrt{p/2}/\|\mathbf{w}_i\|_2^{1-p/2}$, respectively. Then, we can obtain a solution by solving the re-weighted least square minimization problem, i.e.,

$$\min_{\mathbf{W}} \|\mathbf{\Sigma}_1(\mathbf{X}\mathbf{W} - \mathbf{Y})\|_F^2 + \lambda\|\mathbf{\Sigma}_2 \mathbf{W}\|_F^2. \quad (6)$$

Since \mathbf{S}_1 and \mathbf{S}_2 (or $\mathbf{\Sigma}_1$ and $\mathbf{\Sigma}_2$) are functions of \mathbf{W} , we use an iterative algorithm to compute the solution. At each iterative step, \mathbf{S}_1 and \mathbf{S}_2 are fixed first, then \mathbf{W} is obtained according to Eqn. (4), and finally we update \mathbf{S}_1 and \mathbf{S}_2 based on \mathbf{W} . The iterative algorithm is summarized in Algorithm 1, and its convergence is proved in the following subsection.

Algorithm 1. A unified sparse feature selection

1. **Input:** data points $\{\mathbf{x}^i\}_{i=1}^m$ ($\mathbf{x}^i \in \mathbb{R}^n$) and their corresponding label $\{y^i\}_{i=1}^m$; loss function norm order r ; sparse regularization norm order p ; regularization parameter λ ; number of features d to be selected.
 2. Construct \mathbf{X} and \mathbf{Y}
 3. as identity matrices
 4. Set $k = 1$ and initialize $\mathbf{S}_{1_0} \in \mathbb{R}^{m \times m}$ and $\mathbf{S}_{2_0} \in \mathbb{R}^{n \times n}$ as identity matrices
 5. **repeat**
 6. Calculate $\mathbf{W}_k = (\mathbf{X}^T \mathbf{S}_{1_{k-1}} \mathbf{X} + \lambda \mathbf{S}_{2_{k-1}})^{-1} \mathbf{X}^T \mathbf{S}_{1_{k-1}} \mathbf{Y}$
 7. Calculate $\mathbf{E}_k = \mathbf{X}\mathbf{W}_k - \mathbf{Y}$
 8. Update \mathbf{S}_{1_k} , where i -th diagonal elements is $\frac{r/2}{\|e_{i_k}\|_2^{2-r}}$
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9. Update \mathbf{S}_{2_k} , where i -th diagonal elements is $\frac{p/2}{\|\mathbf{w}_{i_k}\|_2^{2-p}}$
 10. Update $k = k + 1$
 11. **Until** convergence
 12. **Output:** Sort all features according to $\|\mathbf{w}_i\|_2$ and select the top largest d features.
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4. Convergence analysis and convergence rate

The objective function $\mathcal{J}(\mathbf{W})$ monotonically decreases at every iteration step and Algorithm 1 finally converges.

Lemma 1. Given any nonzero vectors \mathbf{a} and \mathbf{b} , we have

$$\|\mathbf{b}\|_2^{2\theta} - \theta \cdot \frac{\|\mathbf{b}\|_2^2}{\|\mathbf{a}\|_2^{2-2\theta}} \leq (1 - \theta)\|\mathbf{a}\|_2^{2\theta}, \quad (7)$$

where $0 < \theta < 1$ and the equality holds if and only if $\mathbf{a} = \mathbf{b}$.

Based on Lemma 1 (its proof is available in the supplemental material), we have Lemma 2.

Lemma 2. Given an optimization problem:

$$\min_{\mathbf{Z}} f(\mathbf{Z}) + \|\mathbf{\Sigma} \mathbf{\Phi}(\mathbf{Z})\|_F^2, \quad \text{s.t. } \mathbf{Z} \in \mathcal{F}, \quad (8)$$

where $f(\mathbf{Z})$ is a matrix function of \mathbf{Z} and $\mathbf{\Phi}(\mathbf{Z})$ is a function matrix of \mathbf{Z} , \mathcal{F} is the feasible region. $\mathbf{\Sigma}$ is a diagonal matrix in which the i -th diagonal element is $\sqrt{q/2}/\|\mathbf{\Phi}(\mathbf{Z}_0)_i\|_2^{1-q/2}$ (\mathbf{Z}_0 is one element in \mathcal{F} , $\mathbf{\Phi}(\mathbf{Z}_0)_i$ is the i -th row vector of $\mathbf{\Phi}(\mathbf{Z}_0)$ and $0 < q \leq 2$). If \mathbf{Z}^* is the optimal solution of the above optimization problem Eqn. (8), we have

$$f(\mathbf{Z}^*) + \|\mathbf{\Phi}(\mathbf{Z}^*)\|_{2,q}^q \leq f(\mathbf{Z}_0) + \|\mathbf{\Phi}(\mathbf{Z}_0)\|_{2,q}^q, \quad (9)$$

Proof. Since \mathbf{Z}^* is the optimal solution of Eqn.(8), we have

$$f(\mathbf{Z}^*) + \|\mathbf{\Sigma} \mathbf{\Phi}(\mathbf{Z}^*)\|_F^2 \leq f(\mathbf{Z}_0) + \|\mathbf{\Sigma} \mathbf{\Phi}(\mathbf{Z}_0)\|_F^2. \quad (10)$$

Therefore

$$\begin{aligned} f(\mathbf{Z}^*) + \sum_i \frac{q}{2} \frac{\|\mathbf{\Phi}(\mathbf{Z}^*)_i\|_2^2}{\|\mathbf{\Phi}(\mathbf{Z}_0)_i\|_2^{2-q}} \\ \leq f(\mathbf{Z}_0) + \sum_i \frac{q}{2} \|\mathbf{\Phi}(\mathbf{Z}_0)_i\|_2^q \end{aligned} \quad (11)$$

When $0 < q < 2$, according to Lemma 1, we have

$$\begin{aligned} \sum_i \left(\|\mathbf{\Phi}(\mathbf{Z}^*)_i\|_2^q - \frac{q}{2} \frac{\|\mathbf{\Phi}(\mathbf{Z}^*)_i\|_2^2}{\|\mathbf{\Phi}(\mathbf{Z}_0)_i\|_2^{2-q}} \right) \\ \leq \sum_i \left(1 - \frac{q}{2} \right) \|\mathbf{\Phi}(\mathbf{Z}_0)_i\|_2^q. \end{aligned} \quad (12)$$

Summing Eqn. (11) and Eqn. (12), we obtain

$$f(\mathbf{Z}^*) + \sum_i \|\mathbf{\Phi}(\mathbf{Z}^*)_i\|_2^q \leq f(\mathbf{Z}_0) + \sum_i \|\mathbf{\Phi}(\mathbf{Z}_0)_i\|_2^q. \quad (13)$$

Finally, we obtain

$$f(\mathbf{Z}^*) + \|\Phi(\mathbf{Z}^*)\|_{2,q}^q \leq f(\mathbf{Z}_0) + \|\Phi(\mathbf{Z}_0)\|_{2,q}^q, \quad (14)$$

where the equality holds if and only if $\Phi(\mathbf{Z}^*) = \Phi(\mathbf{Z}_0)$.

When $q = 2$, Σ becomes an identity matrix, the equality in Eqn. (14) still holds. \square

Theorem 1. *The objective function of problem (2) monotonically decreases at every iteration step, i.e.,*

$$\mathcal{J}(\mathbf{W}_k) \leq \mathcal{J}(\mathbf{W}_{k-1}), \quad (15)$$

and it converges to a limit point.

Proof. According to Eqn. (6), we have

$$\begin{aligned} & \|\Sigma_{1_{k-1}}(\mathbf{X}\mathbf{W}_k - \mathbf{Y})\|_F^2 + \lambda \|\Sigma_{2_{k-1}}\mathbf{W}_k\|_F^2 \\ & \leq \|\Sigma_{1_{k-1}}(\mathbf{X}\mathbf{W}_{k-1} - \mathbf{Y})\|_F^2 + \lambda \|\Sigma_{2_{k-1}}\mathbf{W}_{k-1}\|_F^2, \end{aligned} \quad (16)$$

where the i -th diagonal elements of $\Sigma_{1_{k-1}}$ and $\Sigma_{2_{k-1}}$ are $\zeta_{1_{k-1}}^{ii} = \sqrt{r/2} / \|\mathbf{x}_i \mathbf{W}_{k-1} - \mathbf{y}_i\|_2^{1-r/2}$ and $\zeta_{2_{k-1}}^{ii} = \sqrt{p/2} / \|\mathbf{w}_{i_{k-1}}\|_2^{1-p/2}$, respectively.

According to Lemma 2, let $f(\mathbf{W}) = \|\Sigma_{1_{k-1}}(\mathbf{X}\mathbf{W} - \mathbf{Y})\|_F^2$ and $\Phi(\mathbf{W}) = \lambda \mathbf{W}$, we have

$$\begin{aligned} & \|\Sigma_{1_{k-1}}(\mathbf{X}\mathbf{W}_k - \mathbf{Y})\|_F^2 + \lambda \|\mathbf{W}_k\|_{2,p}^p \\ & \leq \|\Sigma_{1_{k-1}}(\mathbf{X}\mathbf{W}_{k-1} - \mathbf{Y})\|_F^2 + \lambda \|\mathbf{W}_{k-1}\|_{2,p}^p. \end{aligned} \quad (17)$$

Then, setting $f(\mathbf{W}) = \lambda \|\mathbf{W}\|_{2,p}^p$ and $\Phi(\mathbf{W}) = \mathbf{X}\mathbf{W} - \mathbf{Y}$, according to Lemma 2 we have

$$\begin{aligned} & \|\mathbf{X}\mathbf{W}_k - \mathbf{Y}\|_{2,r}^r + \lambda \|\mathbf{W}_k\|_{2,p}^p \\ & \leq \|\mathbf{X}\mathbf{W}_{k-1} - \mathbf{Y}\|_{2,r}^r + \lambda \|\mathbf{W}_{k-1}\|_{2,p}^p. \end{aligned} \quad (18)$$

So, $\mathcal{J}(\mathbf{W}_k) \leq \mathcal{J}(\mathbf{W}_{k-1})$, and the equality holds if and only if $\mathbf{W}_k = \mathbf{W}_{k-1}$. Since the lower bound of $\mathcal{J}(\mathbf{W}_k)$ is limited, $\mathcal{J}(\mathbf{W}_k)$ converges to a limit point. \square

Theorem 2. *Sequence $\{\mathbf{W}_k\}$ produced in Algorithm 1 converges, and the limit point is a stationary point of optimization problem (3).*

Proof is presented in the supplemental material.

When $p = 1$ and $r \geq 1$, Eqn. (2) is a convex optimization problem, hence its solution obtained by Algorithm 1 is globally optimal. When $0 < p < 1$ or $0 < r < 1$, it may converge to a local optimum.

The convergence rate of Algorithm 1 is derived as following. If \mathbf{W}^* is the optimal solution of $\min_{\mathbf{W}} \|\mathbf{X}\mathbf{W} - \mathbf{Y}\|_{2,r}^r + \lambda \|\mathbf{W}\|_{2,p}^p$, then the optimal residual $\mathbf{E}^* = \mathbf{X}\mathbf{W}^* - \mathbf{Y}$. When \mathbf{W}^* is sparse, the rows of \mathbf{W}^* can be split into two parts: \mathbf{W}_1^* and \mathbf{W}_2^* , where $\mathbf{W}_2^* = \mathbf{0}$ and \mathbf{W}_1^* is the remainder. In the same way as partitioning \mathbf{W}^* into \mathbf{W}_1^* and \mathbf{W}_2^* , the rows of \mathbf{W}_k and the columns of \mathbf{X} are partitioned into \mathbf{W}_{1_k} and \mathbf{W}_{2_k} , \mathbf{X}_1 and \mathbf{X}_2 , respectively. Similarly, \mathbf{E}^* can be split into \mathbf{E}_1^* and \mathbf{E}_2^* , where $\mathbf{E}_2^* = \mathbf{0}$ and \mathbf{E}_1^* is the remainder, and the rows of \mathbf{E}_k ($\mathbf{E}_k = \mathbf{X}\mathbf{W}_k - \mathbf{Y}$) and the columns of \mathbf{I} ($\in \mathbb{R}^{m \times m}$) are partitioned into \mathbf{E}_{1_k} and \mathbf{E}_{2_k} , \mathbf{L}_1 and \mathbf{L}_2 , accordingly. We define $\mathbf{A}_1 = [\mathbf{X}_1, \mathbf{L}_1]$, $\mathbf{A}_2 = [\mathbf{X}_2, \mathbf{L}_2]$, $\mathbf{U}_{1_k}^T = [\mathbf{W}_{1_k}, -\mathbf{E}_{1_k}]^T$, $\mathbf{U}_{2_k}^T = [\mathbf{W}_{2_k}, -\mathbf{E}_{2_k}]^T$ and $(\mathbf{U}_1^*)^T = [\mathbf{W}_1^*, -\mathbf{E}_1^*]^T$. Then we have Lemma 3.

Lemma 3. *The following inequalities hold in successive iteration steps of Algorithm 1.*

$$\begin{aligned} \|\mathbf{U}_{1_k} - \mathbf{U}_1^*\| & \leq \frac{\frac{2}{p} \|\mathbf{W}_{2_{k-1}}\|_{2,2-p}^{2-p} + \frac{2\lambda}{r} \|\mathbf{E}_{2_{k-1}}\|_{2,2-r}^{2-r}}{\frac{2}{p} \|\mathbf{W}_{1_{k-1}}\|_{2,2-p}^{2-p} + \frac{2\lambda}{r} \|\mathbf{E}_{1_{k-1}}\|_{2,2-p}^{2-p}} \\ & \quad \sqrt{s_0} \|\mathbf{I} - \mathbf{B}^+ \mathbf{B}\| \|\mathbf{A}_1^+ \mathbf{A}_2\|^2 \|\mathbf{U}_1^*\|, \end{aligned} \quad (19)$$

$$\begin{aligned} \|\mathbf{U}_{2_k}\| & \leq \frac{\frac{2}{p} \|\mathbf{W}_{2_{k-1}}\|_{2,2-p}^{2-p} + \frac{2\lambda}{r} \|\mathbf{E}_{2_{k-1}}\|_{2,2-r}^{2-r}}{\frac{2}{p} \|\mathbf{W}_{1_{k-1}}\|_{2,2-p}^{2-p} + \frac{2\lambda}{r} \|\mathbf{E}_{1_{k-1}}\|_{2,2-p}^{2-p}} \\ & \quad \sqrt{s_0} \|\mathbf{I} - \mathbf{B}^+ \mathbf{B}\| \|\mathbf{A}_1^+ \mathbf{A}_2\| \|\mathbf{U}_1^*\|, \end{aligned} \quad (20)$$

where s_0 is the number of columns of \mathbf{A}_1 , and $\mathbf{B} = (\mathbf{I} - \mathbf{A}_1 \mathbf{A}_1^+) \mathbf{A}_2$. (Proof of Lemma 3 is presented in the supplemental material).

According to Lemma 3, we can obtain the convergence order of Algorithm 1.

Theorem 3. *The convergence order of Algorithm 1 is at least $\min(2-p, 2-r)$. (Proof of Theorem 3 is presented in the supplemental material.)*

5. Experiments

5.1 Results based on a synthetic dataset

To investigate how the loss function's parameter r in our method affects the feature selection performance, we generated a synthetic dataset using following procedure. First, we generated n samples with features $\mathbf{X}_1, \mathbf{X}_2 \in \mathbb{R}^{n \times d_1}$, where elements of $\mathbf{X}_1, \mathbf{X}_2$ were randomly generated according to Gaussian distribution $\mathcal{N}(0,1)$. Second, we introduced redundant features $\mathbf{X}_3 = 0.5(\mathbf{X}_1 + \mathbf{X}_2) + \epsilon \in \mathbb{R}^{n \times d_1}$ to the samples, where elements of ϵ were randomly generated according to $\mathcal{N}(0,0.1)$. Third, irrelevant features $\mathbf{X}_4 \in \mathbb{R}^{n \times d_2}$ were injected into the samples, where elements of \mathbf{X}_4 were randomly generated according to uniform distribution $\mathcal{U}(-1,1)$. So, we obtained samples with features $\mathbf{X} = [\mathbf{X}_1, \mathbf{X}_2, \mathbf{X}_3, \mathbf{X}_4] \in \mathbb{R}^{n \times d}$, $d = 3d_1 + d_2$. Then, we generated multi-tasks labels for these samples as $\mathbf{Y}_0 = \mathbf{X}\mathbf{W} + \zeta \in \mathbb{R}^{n \times c}$, where $\mathbf{W} = [\mathbf{W}_1; \mathbf{W}_2; \mathbf{0}] \in \mathbb{R}^{d \times c}$, $\mathbf{W}_1, \mathbf{W}_2 \in \mathbb{R}^{d_1 \times c}$ and their elements were randomly generated according to uniform distribution $\mathcal{U}(0,1)$, and ζ were randomly generated according to $\mathcal{N}(0, 0.5)$. Finally, to simulate outlier samples, we randomly picked a subset of \mathbf{Y}_0 with a percentage of a , and reversed their positive or negative signs, yielding new labels \mathbf{Y}_1 .

Setting $n=2000$, $d_1=100$, $d_2=700$, $c=5$, and $a=0, 0.01$, and 0.1 , we obtained 3 simulated data sets, each of them having 1000 features, among which 200 features were informative.

We evaluated our method using 10-fold cross validation based on the simulated dataset with respect to different $r = 0.5, 1, 2$ by setting $p = 1$. The performance was gauged with

root mean square error (RMSE) between actual values and predicted values base on top 200 selected features. As shown in Figure 2, the least square loss had the best data-fitting performance for samples without outliers. However, it was sensitive to outliers as

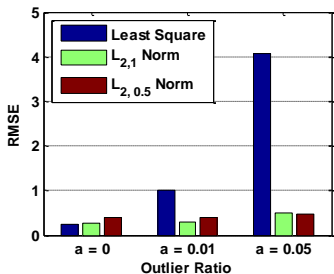


Figure 2: Average RMSE of 10-fold cross-validation for least square loss, $\ell_{2,1}$ -norm loss, and $\ell_{2,0.5}$ -norm loss, respectively.

reflected by relatively larger RMSE when the samples contained 1% and 10% outliers. Not surprisingly, the feature selection with $\ell_{2,r}$ -norm based loss functions ($r \leq 1$) was robust to outliers but might sacrifice data-fitting accuracy. All these results indicate that the loss function’s parameter should be adaptive to the problem under study.

5.2 Classification experiments on real-world datasets

We also evaluated our algorithm, referred to as general sparsity regularized feature selection (GSR-FS), based on 6 publicly available real-world datasets. In particular, 2 datasets were obtained from UCI, including ISOLET and SEMEION. ISOLET is a speech recognition data set with 7797 samples in 26 classes, and each sample has 617 features. SEMEION contains 1593 handwritten images from ~80 persons, stretched in a rectangular box of 16×16 . Three face image datasets were obtained from AR, ORL, and the frontal pose sub-dataset (09) of CMU-PIE. Particularly, AR has 1680 samples with 2000 features, ORL contains 400 samples with 92×112 pixels as features, and the CMU-PIE subset contains images of 64 persons with different illuminations. The last dataset contains confusable hand writing images 4 and 9, obtained from MNIST.

We compared our method with 4 sparsity regularized feature selection algorithms, including MTFS (Argyriou &

Evgeniou, 2007; Liu, et al., 2009; Obozinski, et al., 2006), RFS (Nie, et al., 2010; Xiang, et al., 2012), an extended RFS (E-RFS) (Wang & Chen, 2013), and RK1U (Zhang, et al., 2014). We also compared our method with two filter feature selection methods, namely ReliefF (Kira & Rendell, 1992) and mRMR (Peng, et al., 2005).

In our experiments, we first normalized all the features to have 0 mean and unit standard deviation. Then, 10 trials were carried out on each dataset for feature selection. In each trial, each dataset was randomly split into training and testing subsets with a ratio of 6:4. Classification accuracy was used to evaluate the feature selection methods. Particularly, linear SVM (Chang & Lin, 2011) was used to build classifiers based on the selected features. The parameter C of linear SVM classifiers was tuned using a cross-validation strategy by searching a candidate set of [10-3, 10-2, 10-1, 1, 101, 102]. The regularized parameter λ in our algorithm, MTFS, RFS and RK1U was tuned using the same cross-validation strategy by searching a candidate set of [10-3, 10-2, 10-1, 1, 101, 102].

Our algorithm has 2 hyper parameters r and p . For evaluating the impact of p on the sparsity and directly comparing our method with MTFS, RFS and RK1U, we evaluated our algorithm by setting $p = 1, 0.75, 0.5$ and 0.25 . Since the loss function with smaller r is more robust to outliers but a larger r of the loss function may yield better data-fitting performance, in our experiments r was tuned by cross-validation with a candidate set of [0.5, 1, 2]. For the E-RFS and RK1U, $p = 0.5$ since it had better classification performance than other values (Wang & Chen, 2013; Zhang, et al., 2014).

Table 2 and Figure 3 summarize the average classification performance of classifiers built on features selected by different methods in 10 trials. The average classification accuracy rates with top [10, 20, ..., 100] features are shown in Figure 3. Table 2 summarizes mean and standard deviation of the classification rates in 10 trails for classifiers built on the top 50 features. These results demonstrated that our method with different p achieved overall the best classification accuracy on most of the datasets, especially when $p = 0.75, 0.5$. When of the flexibility in our data-loss function. When $p = 0.5$, our method performed better than E-RFS and RK1U. Not surprisingly, the sparsity regularization methods had better performance than filter methods.

Table 2: Mean and standard deviation of the classification accuracy (% , mean \pm std) of Linear-SVM classifiers built on the top 50 features selected by different algorithms on different datasets.

Algorithm	ReliefF	mRMR	MTFS	RFS	Extended RFS (p=0.5)	RK1U (p=0.5)	GSR-FS (p=1.0)	GSR-FS (p=0.75)	GSR-FS (p=0.5)	GSR-FS (p=0.25)
ISOLET	77.02 \pm 0.82	85.10 \pm 0.62	90.40 \pm 0.68	91.38 \pm 0.73	91.36 \pm 0.56	92.50 \pm 0.79	92.98 \pm 0.44	93.92 \pm 0.46	94.10\pm0.29	93.93 \pm 0.40
SEMEION	78.70 \pm 1.39	78.75 \pm 1.82	84.11 \pm 1.12	85.82 \pm 1.25	85.89 \pm 1.61	85.45 \pm 1.01	86.24 \pm 1.45	86.97\pm1.13	86.87 \pm 0.89	86.58 \pm 1.17
AR	57.89 \pm 4.58	87.77 \pm 1.36	85.24 \pm 2.99	88.90 \pm 1.41	88.36 \pm 1.82	87.90 \pm 1.88	92.38 \pm 1.24	94.49 \pm 0.56	94.51\pm0.96	94.06 \pm 0.89
ORL	56.50 \pm 6.12	87.81 \pm 3.23	74.00 \pm 3.60	80.25 \pm 2.74	88.44 \pm 2.11	90.13 \pm 1.81	90.88 \pm 2.83	91.50\pm2.27	91.13 \pm 2.93	90.06 \pm 2.33
CMU-PIE	75.35 \pm 1.50	89.91 \pm 1.04	85.41 \pm 1.39	90.83 \pm 0.79	90.98 \pm 1.01	91.70 \pm 0.97	93.01 \pm 0.87	93.85\pm0.83	93.67 \pm 1.08	93.39 \pm 0.79
MNIST	90.62 \pm 0.35	92.55 \pm 0.19	95.30 \pm 0.21	94.25 \pm 0.38	94.04 \pm 0.55	95.31 \pm 0.20	95.53 \pm 0.21	95.85\pm0.24	95.53 \pm 0.18	95.49 \pm 0.35

Table 3: Running Time (unit: second) Taken by different Algorithms

Algorithm	Relieff	mRMR	RFS	Extended RFS(p=0.5)	RK1U (p=0.5)	GSR-FS (p=1.0)	GSR-FS (p=0.75)	GSR-FS (p=0.5)	GSR-FS (p=0.25)
ISOLET	319.48	52.85	2954.07	2100.20	6467.20	26.70	24.70	21.47	17.01
SEMEION	6.18	4.70	56.38	32.93	101.69	11.11	5.52	5.58	5.60
AR	37.13	24.67	67.44	48.92	2070.54	62.08	30.19	26.36	20.49
ORL	23.04	157.69	12.01	7.81	5750.64	12.11	11.97	4.60	3.21
CMU-PIE	89.35	62.24	66.47	67.31	4563.38	49.38	53.22	35.64	19.43
MNIST	676.21	48.83	9275.84	6199.82	1050.98	26.87	22.56	18.81	18.87

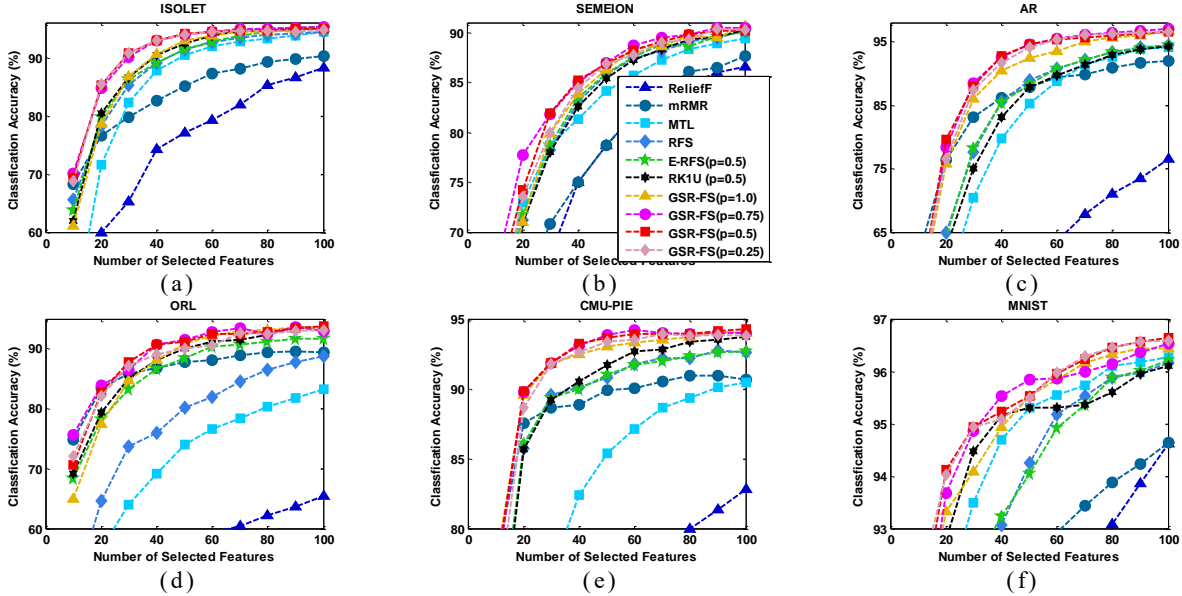


Figure 3: Average classification accuracy of 10 trials for classifiers built on the selected features by different algorithms. The results shown were obtained on (a) ISOLET, (b)SEMEION, (c) AR, (d) ORL, (e) CMU-PIE, and (f) MNIST.

5.3 Computational cost

We also compared our algorithm with other methods with respect to their computation cost². The convergence of all the sparse feature selection algorithms was determined based on the same criterion: the change of objective function value is less than 10^{-4} between 2 successive iteration steps with the regularized parameter $\lambda=1$. And the filter algorithms run until top 100 features were selected. The parameter r in our algorithm was set to 2. We run different methods on a desktop with a Intel i7-4470 cpu, 3.4GHz and 8G RAM. The computation costs of different algorithms are summarized in Table 3. As shown in Table 3, our algorithm was faster than other sparse feature selection algorithms on most of the datasets, and had similar costs as mRMR and ReliefF. Particularly, RK1U and our method achieved similar classification performance on some datasets, but the computational time of RK1U is more than 50 times longer than ours on average.

6. Discussions and Conclusion

We have presented a general framework for sparsity regularized feature selection and a novel iterative reweighted least square minimization optimization algorithm. Several existing sparsity regularized feature selection methods could be treated as its special cases. The objective function of our method consists of a $\ell_{2,r}$ -norm ($0 < r \leq 2$) based loss function and a $\ell_{2,p}$ -norm ($0 < p \leq 1$) sparse regularization, resulting in an adaptive setting for handling outliers by turning its parameters. Such flexibility could improve feature selection performance as demonstrated by the experimental results. The novel IRLS algorithm is capable of solving problems with multiple non-smooth functions, and could find its applications in other fields. We will extend our method to constrained optimization problems and investigate how to choose optimal parameters r and p besides cross-validation.

² MTFS was implemented in C, and other algorithms were implemented in Matlab. So, we did not directly compare our algorithm with MTFS. However, RK1U was faster than MTFS (M. Zhang, et al, 2014), and our method is faster than RK1U.

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