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# The Support Vector Regression with Adaptive Norms

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#### **Abstract**

This study proposes a new method for regression  $-l_p$ -norm support vector regression ( $l_p$  SVR). Some classical SVRs minimize the hinge loss function subject to the  $l_2$ -norm or  $l_1$ -norm penalty. These methods are non-adaptive since their penalty forms are fixed and pre-determined for any types of data. Our new model is an adaptive learning procedure with lp-norm (0 < p < 1), where the best p is automatically chosen by data. By adjusting the parameter p,  $l_p$  SVR can not only select relevant features but also improve the regression accuracy. An iterative algorithm is suggested to solve the  $l_p$  SVR efficiently. Simulations and real data applications support the effectiveness of the proposed procedure.

Keywords: Regression; Support vector machine; Norm; Feature selection;

### 1. Introduction

Support vector machines (SVMs), being computationally powerful tools for pattern classification and regression, have been successfully applied to a variety of real-world problems([1]- [6]). Regards to the support vector regression (SVR), some classical SVRs minimize the hinge loss function subject to the  $l_2$ -norm or  $l_1$ -norm penalty ([7]). We call them  $l_2$  SVR or  $l_1$  SVR correspondingly. These methods are non-adaptive since their penalty forms are fixed and pre-determined for any types of data.

Recently,  $l_p$ -norm ( $p \in (0,1)$ ) attracts great attention in the optimization framework, the idea that using  $l_p$ -norm can find sparse solutions is considered in [8]-[11]. Correspondingly, [12]-[17] propose  $l_p$ -norm ( $0 ) support vector machine for classification (<math>l_p$  SVC), which replace the  $l_2$ -norm penalty by the  $l_p$ -norm ( $p \in (0,1)$ ) penalty in the objective function in the primal problem in the standard linear  $l_2$  SVC. Compared with SVC with a fixed norm,  $l_p$  SVC is desired for feature selection since it can automatically select relevant features by adjusting the parameter p. However,  $l_p$  SVC is used only to solve classification problems. This motivates us to consider a new model with  $l_p$ -norm for regression problems.

This paper proposes a new method for regression  $-l_p$ -norm support vector regression  $(l_p \text{ SVR})$ , which replaces  $l_2$ -norm by  $l_p$ -norm  $(0 in the classical <math>l_2 \text{ SVR}$ . Our new model is an adaptive learning procedure with  $l_p$ -norm (0 , where the best <math>p is automatically chosen by data. By adjusting the parameter p,  $l_p \text{ SVR}$  can not only select relevant features but also improve the regression accuracy. In order to solve the non-convex problem in our model, an efficient algorithm is constructed using the successive linear approximation algorithm (SLA)([18]).

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Now we describe our notation. All vectors are column vectors unless transposed to a row vector by a super script T. For a vector x in  $R^n$ ,  $[x]_i(i=1,2,\cdots,n)$  denotes the i-th component of x. |x| denotes a vector in  $R^n$  of absolute value of the components of x.  $||x||_p$  denotes that  $(|[x]_1|^p + \cdots + |[x]_n|^p)^{\frac{1}{p}}$ . Strictly speaking,  $||x||_p$  is not a general norm when  $0 , but we still follow this term <math>l_p$ -norm, because the forms are same except that the values of p are different.  $||x||_0$  is the number of nonzero components of x. For two vectors  $x \in R^n$  and  $y \in R^n$ ,  $(x \cdot y)$  denotes the inner product of x and y.

This paper is organized as follows. In section 2, the  $l_p$  SVR is introduced. In section 3, the SLA is proposed to solve  $l_p$  SVR. In section 4, the lower bounds for the absolute value of nonzero entries in any local optimal solution is established. In section 5, numerical experiments are given to demonstrate the effectiveness of our method. We conclude this paper in section 6.

## 2. $l_p$ Support Vector Regression

Suppose that the training set *T* is given by

$$T = \{(x_1, y_1), \dots, (x_l, y_l)\} \in (R^n \times R)^l,$$
(1)

where  $x_j \in R^n$ ,  $y_j \in R$   $(j = 1, \dots, l)$ , the linear regression problem is to find a decision function  $f(x) = (w \cdot x) + b$  to derive the value of y for any x by the function y = f(x).

In the classical  $l_2$  SVR, the decision function is decided by the solution to the following optimization problem:

$$\min_{w,b,\xi,\eta} \frac{1}{2} ||w||_2^2 + C \sum_{i=1}^l (\eta_i + \xi_i), \qquad (2)$$

s.t. 
$$((w \cdot x_i) + b) - y_i \le \varepsilon + \eta_i , i = 1, \dots, l ,$$
 (3)

$$y_i - ((w \cdot x_i) + b) \le \varepsilon + \xi_i, i = 1, \dots, l,$$
(4)

$$\xi_i, \eta_i \ge 0 , i = 1, \cdots, l. \tag{5}$$

Replacing the first term  $\frac{1}{2}||w||_2^2$  in the objective function of the above problem by  $||w||_p^p$  (0 < p < 1),  $l_p$  SVR proposes the following problem:

$$\min_{w,b,\xi,\eta} ||w||_p^p + C \sum_{i=1}^l (\eta_i + \xi_i),$$
 (6)

s.t. 
$$((w \cdot x_i) + b) - y_i \le \varepsilon + \eta_i, \quad i = 1, \dots, l,$$
 (7)

$$y_i - ((w \cdot x_i) + b) \le \varepsilon + \xi_i, i = 1, \dots, l,$$
(8)

$$\xi_i, \eta_i \ge 0 , i = 1, \cdots, l, \tag{9}$$

where C (C > 0), p ( $0 ) and <math>\varepsilon$  ( $\varepsilon > 0$ ) are parameters. The algorithm of  $l_p$  SVR is described as follows:

### Algorithm 1

- 1. Give the training set  $T = \{(x_1, y_1), \dots, (x_l, y_l)\} \in (R^n \times R)^l$ , where  $x_i \in R^n, y_i \in R, i = 1, \dots, l$ ;
- 2. Select proper parameters C, p,  $\varepsilon$ , where C > 0,  $0 and <math>\varepsilon > 0$ ;
- 3. Solve problem (6)-(9) and get the solution  $(w^*, b^*)$ ;
- 4. Select the feature set:  $\{i|[w^*]_i \neq 0, (i = 1, \dots, n)\};$
- 5. Construct the decision function  $y = (\tilde{w}^* \cdot x) + b^*$ , where the components of  $\tilde{w}^*$  are nonzero components of  $w^*$  and the components of  $\tilde{x}$  are also corresponding to nonzero components of  $w^*$ .

## 3. The SLA for problem (6)-(9)

Consider the problem (6)-(9), the objective function is not differentiable, because of the absolute value in the first item. In order to make this problem smooth, we introduce the variable  $\nu = ([\nu]_1, \dots, [\nu]_n)^T$ , and get the

following equivalent problem:

$$\min_{w,v,b,\xi,\eta} \qquad \sum_{i=1}^{n} [v]_{i}^{p} + C \sum_{i=1}^{l} (\eta_{i} + \xi_{i}), \qquad (10)$$

s.t. 
$$((w \cdot x_i) + b) - y_i \le \varepsilon + \eta_i, \quad i = 1, \dots, l,$$
 (11)

$$y_i - ((w \cdot x_i) + b) \le \varepsilon + \xi_i, i = 1, \dots, l, \tag{12}$$

$$\xi_i, \eta_i \ge 0 , i = 1, \dots, l, \tag{13}$$

$$-v < w < v. \tag{14}$$

When  $p \in (0, 1)$ , we note that the problem (10)-(14) is differentiable, but not convex. In fact, it is the minimization of a concave objective function over a polyhedral set. Even though it is difficult to find a global solution to this problem, a fast successive linear approximation (SLA) algorithm ([18]) terminates finitely at a stationary point which satisfies the necessary optimality condition for problem (10)-(14). For convenience we state the SLA algorithm below.

## Algorithm 2 (SLA for problem (10)-(14))

- 1. Select the proper parameters C > 0, 0 0 and a precision  $\delta(0 < \delta << 1)$ , start with a random  $v^0 = ([v]_1^0, [v]_2^0, \cdots, [v]_n^0)^T$  and let k = 1; 2. Solve the following problem

$$\min_{w,v,b,\xi,\eta} \qquad \sum_{i=1}^{n} [v^{k-1}]_{i}^{p-1}[v]_{i} + C \sum_{i=1}^{l} (\eta_{i} + \xi_{i}) , \qquad (15)$$

s.t. 
$$((w \cdot x_i) + b) - y_i \le \varepsilon + \eta_i, \quad i = 1, \dots, l,$$
 (16)

$$y_i - ((w \cdot x_i) + b) \le \varepsilon + \xi_i, i = 1, \dots, l,$$

$$(17)$$

$$\xi_i, \eta_i \ge 0 , i = 1, \dots, l, \tag{18}$$

$$-\nu \le w \le \nu. \tag{19}$$

 $-\nu \le w \le \nu$ . where  $(\nu^{k-1})^{p-1} = ([\nu^{k-1}]_1^{p-1}, \dots, [\nu^{k-1}]_n^{p-1})^T$ , and get its solution  $(w^k, b^k, \eta^k, \xi^k, \nu^k)$ ;

3. If 
$$\left| \sum_{i=1}^{n} [v^{k-1}]_i^{p-1} ([v^k]_i - [v^{k-1}]_i^{p-1}) + C \sum_{i=1}^{l} (\eta_i^k - \eta_i^{k-1} + \xi_i^k - \xi_i^{k-1}) \right| < \delta(0 < \delta \ll 1)$$
, then stop and get the solution  $w^* = w^k$ ,  $b^* = b^k$ ; Otherwise, let  $k = k + 1$  and go back to step 2.

## 4. The lower bounds for nonzero components in solutions

In Algorithm 1, it is easy to see that our  $l_p$  SVR can accomplish feature selection and regression simultaneously. Feature selection needs to find the nonzero components of the solution to the problem (6)-(9). However, usually the above Algorithm 2 can only provide an approximate local solution where nonzero components in the solution can not be identified theoretically. Using a similar strategy in [8], we get the following theorem 1, which can be used to identify nonzero components in any local optimal solutions to the problem (6)-(9), even though the Algorithm 2 can only find the approximate local optimal solution.

**Theorem 1** For any local optimal solution  $(w^*, b^*, \xi^*)$  to the problem (6)-(9), we have

$$|[w^*]_j| \ge (p/(C\sum_{i=1}^l |x_i|_j))^{\frac{1}{1-p}}, j = 1, 2, \dots, n.$$

**Proof.** Suppose  $||w^*||_0 = k$ ,  $(1 < k \le n)$ , without loss of generality, let  $w^* = ([w]_1^*, \dots, [w]_k^*, 0, \dots, 0)^T$  where  $[w]_i^* \ne 0$ ,  $i = 1, \dots, k$ . Let  $\tilde{w}^* = ([w]_1^*, \dots, [w]_k^*)^T$ ,  $\tilde{x}_i = ([x_i]_1, \dots, [x_i]_k)^T \in R^k$ ,  $i = 1, \dots, l$ , we consider a new optimization problem:

$$\min_{\tilde{w}, \tilde{b}, \tilde{\xi}, \tilde{\eta}} \qquad \|\tilde{w}\|_p^p + C \sum_{i=1}^l (\tilde{\eta}_i + \tilde{\xi}_i) , \qquad (20)$$

s.t. 
$$((\tilde{w} \cdot \tilde{x}_i) + \tilde{b}) - y_i \le \varepsilon + \tilde{\eta}_i, \quad i = 1, \dots, l,$$
 (21)

$$y_i - ((\tilde{w} \cdot \tilde{x}_i) + b) \le \varepsilon + \tilde{\xi}_i, i = 1, \dots, l,$$
 (22)

$$\tilde{\xi}_i, \tilde{\eta}_i \ge 0$$
,  $i = 1, \dots, l$ , (23)

where  $\tilde{w} \in R^k, \tilde{b} \in R, \tilde{\eta} \in R^l, \tilde{\xi} \in R^l$ . Obviously,  $(\tilde{w}^*, b^*, \eta^*, \xi^*)$  is a local minimizer of problem (20)-(23). According to the KKT condition, there exist Lagrange multipliers  $\alpha_i^*, \beta_i^* (i = 1, \dots, l)$  satisfy:

$$p(|\tilde{w}^*|^{p-1} \cdot sign(\tilde{w}^*)) - \sum_{i=1}^{l} (\beta_i^* - \alpha_i^*) \tilde{x}_i = 0,$$
(24)

$$0 \le \alpha_i^* \le C, \ 0 \le \beta_i^* \le C. \tag{25}$$

According to (24), we have

$$p(|\tilde{w}^*|^{p-1} \cdot sign(\tilde{w}^*)) = \sum_{i=1}^{l} (\beta_i - \alpha_i)\tilde{x}_i.$$

Furthermore, by (25), we have

$$p \, |\tilde{w}^*|^{p-1} = \left| \sum_{i=1}^l (\beta_i - \alpha_i) \tilde{x}_i \right| \leq \sum_{i=1}^l |\beta_i - \alpha_i| \, |\tilde{x}_i| \leq C \sum_{i=1}^l |x_i|, 0$$

So, 
$$\left| w_j^* \right| \ge (p/(C\sum_{i=1}^l |x_i|_j))^{\frac{1}{1-p}}$$
, for  $j = 1, \dots, n$ .

According to Theorem 1, we can identify the nonzero components of the local optimal solution to (6)-(9). Based on the Algorithm 2 and Theorem 1, the new algorithm is established as follows:

## Algorithm 3 ( $l_p$ -SVR):

- 1. Give the training set  $T = \{(x_1, y_1), \dots, (x_l, y_l)\} \in (R^n \times R)^l$ , where  $x_i \in R^n, y_i \in R, i = 1, \dots, l$ ;
- 2. Select proper parameters C, p,  $\epsilon$ , where C > 0, 0 ;
- 3. Solve problem (6)-(9) by Algorithm 2 and get the solution  $(w^*, b^*)$ ;
- 4. Compute  $L_j = (p/(C\sum_{i=1}^{l} |x_i|_j))^{\frac{1}{1-p}}, j = 1, \dots, n$ ; select the feature index set:  $F' = \{i | [w^*]_i | \ge L_i, i = 1, \dots, n\}$ ;
- 5. Construct the decision function  $f(x) = \text{sgn}((\tilde{w}^* \cdot \tilde{x}) + b^*)$ , where  $\tilde{w}^*$  are composed by the components in the F' of  $w^*$  and the components of  $\tilde{x}$  are also corresponding to components in the feature set F' of  $w^*$ .

In the following section, our experiments are conducted according to the algorithm 3.

### 5. Numerical experiments

In this section, some experiments on simulation datasets and real datasets are conducted respectively, by comparing  $l_p$  SVR with  $l_2$  SVR,  $l_1$  SVR. Note that, the performance of each method depend on the parameters (C, p and  $\varepsilon$  in  $l_p$  SVR; C and  $\varepsilon$  in  $l_2$  SVR and  $l_1$  SVR). Therefore, these parameters should be adjusted properly. In our experiments, the best value of these parameters are chosen by five-fold cross validation. C is obtained through searching in the range  $2^{-7} - 2^4$ , p is chosen from 0.1 - 0.9 and  $\varepsilon$  is chosen from 0.01 to 0.1.

In order to evaluate the performance of algorithms, some evaluation criteria ([19], [20]) commonly used should be introduced in the following:

$$MSE = \frac{1}{m} \sum_{i=1}^{m} (f(x_i) - y_i)^2, MAPE = \frac{\sum_{i=1}^{m} \frac{|y_i - f(x_i)|}{|y_i|}}{m} \times 100\%, NMSE = \frac{\sum_{i=1}^{m} (y_i - f(x_i))^2}{\sum_{i=1}^{m} (y_i - \bar{y})^2},$$

$$R^{2} = \frac{\sum_{i=1}^{m} (f(x_{i}) - \bar{y})^{2}}{\sum_{i=1}^{m} (y_{i} - \bar{y})^{2}}, r = \frac{m \sum_{i=1}^{m} f(x_{i}) y_{i} - \sum_{i=1}^{m} f(x_{i}) \sum_{i=1}^{m} y_{i}}{\sqrt{(m \sum_{i=1}^{m} f(x_{i})^{2} - (\sum_{i=1}^{m} f(x_{i}))^{2})(m \sum_{i=1}^{m} y_{i}^{2} - (\sum_{i=1}^{m} y_{i})^{2})}},$$

where m is the number of testing samples,  $f(x_i)$  denotes the predict value of  $x_i$  and  $\bar{y}$  is the average value of  $y_1, \dots, y_l$ .

#### 5.1. Simulation datasets

The simulation datasets are generated as follows. The inputs  $x_i \in \mathbb{R}^n$  are stochastic vectors independently generated in [0, 1], and the numbers of samples and features are described in Table 1. The outputs are determined

Table 1. Description of simulation datasets

Datasets	No. of samples	No. of features
1	400	50
2	500	60
3	500	50
4	100	120

by some simple functions. For example, in dataset 1, the output  $y_i$  is given by

$$y_i = 2[x_i]_1 + 3[x_i]_2 + 4[x_i]_3 + 0.1 \times rand(1);$$

In dataset 2.

$$y_i = 8[x_i]_1 - 7[x_i]_5 + 6[x_i]_9 - 5[x_i]_{13} + 4[x_i]_{20} - 3[x_i]_{31} + 2[x_i]_{45} - [x_i]_{49} + rand(1);$$

In dataset 3,

$$y_i = 100[x_i]_3 + 20[x_i]_{17} + 3[x_i]_{21} + 0.4[x_i]_{36} + 0.05[x_i]_{44};$$

In dataset 4,

$$y_i = 2[x_i]_1 + 3[x_i]_2 + 4[x_i]_3 + rand(1).$$

The results on four datasets are illustrated in Table 2. We show the effectiveness of  $l_p$  SVR from two aspects: feature selection and regression accuracy. On the one hand, from the data in 3th column, it is easy to see that  $l_p$  SVR selects the minimal features. On the other hand, the data in 4th-8th column show that  $l_p$  SVR derives the smallest MSE, MAPE, NMSE, and the largest  $R^2$ , r among these methods in most datasets. This indicates that the statistical information in these datasets is well presented by our  $l_p$  SVR with fairly small feature sets and regression errors.

Table 2. Results on simulation datasets

Datasets	Regressor	No. of selected	MSE	MAPE	NMSE	$R^2$	r
		features					
1	$l_p$ SVR	3	0.0008	0.0064	0.0003	1.0080	0.9999
	$\hat{l_2}$ SVR	50	0.0017	0.0084	0.0006	1.0059	0.9997
	$l_1$ SVR	3	0.0008	0.0062	0.0003	0.9963	0.9999
2	$l_p$ SVR	8	0.0811	0.1561	0.0044	0.9848	0.9978
	$\hat{l_2}$ SVR	60	0.0871	0.1623	0.0048	0.9999	0.9976
	$l_1$ SVR	16	63.0302	5.6449	3.4472	3.6134	0.7724
3	$l_p$ SVR	5	0.0763	0.0346	0.0221	0.9596	0.9889
	$l_2$ SVR	50	0.1038	0.0393	0.0301	0.9622	0.9849
	$l_1$ SVR	16	0.3486	0.0802	0.1011	1.0428	0.9879
4	$l_p$ SVR	3	0.1222	0.0780	0.0377	0.9023	0.9843
	$\hat{l_2}$ SVR	120	1.2196	0.2420	0.3758	0.6735	0.7999
	$l_1$ SVR	3	0.1420	0.0734	0.0438	0.7510	0.9875

#### 5.2. Real datasets

For further evaluation of our method, we choose four real datasets: "bodyfat", "cpusmall", "housing" and "insurance", which are commonly used in testing machine learning algorithms. More detailed description can be found in Table 3.

Table 4 lists the results of three methods on four real datasets. It can be seen that our  $l_p$  SVR can accomplish the desired feature selection and achieve the good regression accuracy. The reason maybe that it can balance these two aspects better than the other two methods by adjusting the parameter p.

Table 3. Description of real datasets

Datasets	No. of samples	No. of features
bodyfat	252	14
cpusmall	500	12
housing	452	13
insurance	500	85

Table 4. Results on real datasets

Datasets	Regressor	No. of selected	MSE	MAPE	NMSE	$R^2$	r
		features					
bodyfat	$l_p$ SVR	1	0.0000	0.0017	0.0004	1.0017	0.9998
	$l_2$ SVR	14	0.0000	0.0016	0.0004	1.0041	0.9998
	$l_1$ SVR	1	0.0305	0.0981	1.1352	0.0550	0.1710
cpusmall	$l_p$ SVR	8	0.0010	0.0124	0.1508	0.7845	0.9354
	$l_2$ SVR	12	0.0009	0.0122	0.1445	0.7973	0.9378
	$l_1$ SVR	2	0.0992	0.1625	15.4089	14.7173	0.2143
housing	$l_p$ SVR	3	0.0097	0.0446	0.2607	0.7223	0.8607
	$\hat{l_2}$ SVR	13	0.0097	0.0416	0.2587	0.7710	0.8641
	$l_1$ SVR	3	0.1372	0.2673	3.6738	3.7220	0.7769
insurance	$l_p$ SVR	9	2.3464	0.1123	0.0141	0.9411	0.9933
	$l_2$ SVR	83	2.8524	0.1121	0.0172	0.9916	0.9917
	$l_1$ SVR	2	3.3524	0.1556	0.0198	0.9070	0.9927

### 6. Conclusions

For regression problems, a new model  $l_p$  SVR is proposed in this paper. The main contribution is that the desired feature selection and good regression performance are implemented simultaneously by introducing the adaptive norms –  $l_p$ -norm, where the parameter p can be chosen flexibly in (0,1) by data. Computational comparisons between our  $l_p$  SVR and other popular methods including  $l_2$  SVR and  $l_1$  SVR indicate the effectiveness of our method. We believe that its good performance mainly comes from the fact that the parameter p is adjusted properly.

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