On the Convergence of the Decomposition Method for Support Vector Machines

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Abstract

The decomposition method is currently one of the major methods for solving support vector machines (SVM). Its convergence properties have not been fully understood. The general asymptotic convergence was first proposed by Chang et al. [3]. However, their working set selection does not coincide with existing implementation. A later breakthrough by Keerthi and Gilbert [12] proved the finite termination for practical cases while the size of the working set is restricted to two. In this paper, we prove the asymptotic convergence of the algorithm used by the software SVMlight [11] and other later implementation. The size of the working set can be any even number. Extensions to other SVM formulations are also discussed.

Keywords
Support vector machines, decomposition methods, classification.

I. Introduction

The support vector machine (SVM) is a new and promising technique for classification. Surveys of SVM are, for example, by Vapnik [27], [28] and Schölkopf et al. [23]. Given training vectors \( x_i \in \mathbb{R}^n, i = 1, \ldots, l \), in two classes, and a vector \( y \in \mathbb{R}^l \) such that \( y_i \in \{1, -1\} \), the support vector technique requires the solution of the following optimization problem:

\[
\min \frac{1}{2} \alpha^T Q \alpha \quad - e^T \alpha \\
0 \leq \alpha_i \leq C, i = 1, \ldots, l, \\
y^T \alpha = 0,
\]  

(1)
where $e$ is the vector of all ones, $C$ is the upper bound of all variables, and $Q$ is an $l$ by $l$ positive semidefinite matrix. Training vectors $x_i$ are mapped into a higher (maybe infinite) dimensional space by the function $\phi$ and $Q_{ij} \equiv y_i y_j K(x_i, x_j)$ where $K(x_i, x_j) \equiv \phi(x_i)^T \phi(x_j)$ is the kernel.

The difficulty of solving (1) is the density of $Q$ because $Q_{ij}$ is in general not zero. In this case, $Q$ becomes a fully dense matrix so a prohibitive amount of memory is required to store the matrix. Thus traditional optimization algorithms such as Newton, Quasi Newton, etc., cannot be directly applied. Several authors (for example, Osuna et al. [18], Joachims [11], Platt [19], and Saunders et al. [22]) have proposed decomposition methods to conquer this difficulty. The basic concept of this method is as follows:

**Algorithm I.1 (Decomposition method)**

1. Given a number $q \leq l$ as the size of the working set. Find $\alpha^1$ as the initial solution. Set $k = 1$.
2. If $\alpha^k$ is an optimal solution of (1), stop. Otherwise, find a working set $B \subset \{1, \ldots, l\}$ whose size is $q$. Define $N \equiv \{1, \ldots, l\} \setminus B$ and $\alpha_B^k$ and $\alpha_N^k$ to be sub-vectors of $\alpha^k$ corresponding to $B$ and $N$, respectively.
3. Solve the following sub-problem with the variable $\alpha_B$:

   $$\min \frac{1}{2} \alpha_B^T Q_{BB} \alpha_B - (e_B - Q_{BN} \alpha_N^k)^T \alpha_B$$

   $$0 \leq (\alpha_B)_i \leq C, i = 1, \ldots, q,$$

   $$y_B^T \alpha_B = -y_N^T \alpha_N^k.$$  \hspace{1cm} (2)

   where $[Q_{BB} \; Q_{BN} \; Q_{NN}]$ is a permutation of the matrix $Q$.
4. Set $\alpha_B^{k+1}$ to be the optimal solution of (2) and $\alpha_N^{k+1} \equiv \alpha_N^k$. Set $k \leftarrow k + 1$ and goto Step 2.

The basic idea of the decomposition method is that in each iteration, the indices $\{1, \ldots, l\}$ of the training set are separated to two sets $B$ and $N$, where $B$ is the working set and $N = \{1, \ldots, l\} \setminus B$. The vector $\alpha_N$ is fixed so the objective value becomes $\frac{1}{2} \alpha_B^T Q_{BB} \alpha_B - (e_B - Q_{BN} \alpha_N)T \alpha_B + \frac{1}{2} \alpha_N^T Q_{NN} \alpha_N - e_N^T \alpha_N$. Then a sub-problem with the variable $\alpha_B$, i.e. (2), is solved. Note that $B$ is updated in each iteration. To simplify the notation, we simply use $B$ instead of $B^k$. 


An important issue of the decomposition method is to select the working set $B$ in each iteration (Step 2 of Algorithm I.1). Among existing methods, Osuna et al. [18], and Saunders et al. [22] find the working set by choosing elements which violate the Karush-Kuhn-Tucker (KKT) condition. Platt’s Sequential Minimal Optimization (SMO) [19] restricts the size of the working set to be two. The advantage is that (2) becomes a small problem so no optimization software is required in practice. His working selection includes some heuristics. A systematic way is proposed by Joachims [11] where he restricts $q$ to be an even number. In his software $SVM^{light}$, the following problem with the variable $d$ is solved:

$$
\begin{align}
\min & \quad \nabla f(\alpha^k)^T d \\
& \quad y^T d = 0, \quad -1 \leq d_i \leq 1, i = 1, \ldots, l, \\
& \quad d_i \geq 0, \text{ if } (\alpha^k)_i = 0, \quad d_i \leq 0, \text{ if } (\alpha^k)_i = C, \\
& \quad |\{d_i \mid d_i \neq 0\}| \leq q,
\end{align}
$$

(3a, 3b, 3c)

where we represent $f(\alpha) \equiv \frac{1}{2} \alpha^T Q \alpha - e^T \alpha$, $\alpha^k$ is the solution at the $k$th iteration, and $\nabla f(\alpha^k)$ is the gradient of $f(\alpha)$ at $\alpha^k$. Note that $|\{d_i \mid d_i \neq 0\}|$ means the number of components of $d$ which are not zero. The constraint (3c) implies that a direction $d$ involving only $q$ variables is obtained. Then components of $\alpha^k$ with non-zero $d_i$ are included in the working set $B$ which is used to construct the sub-problem (2). Note that $d$ is only used for identifying $B$ but not as a search direction. In Joachims’ original paper, $|\{d_i \mid d_i \neq 0\}| = q$ instead of (3c) was used. Thus practically the decomposition method always picks $q$ elements in each iteration. It was first pointed out in [3] that in theory $q$ nonzero elements may not be always available so an inequality (3c) was proposed.

Joachims [11] used the following procedure to solve (3):

**Algorithm I.2 (SVM$^{light}$’s working set selection)**

1. Sort $y_i \nabla f(\alpha^k)_i$ in the decreasing order.
2. From the top of the sorted list sequentially set $d_i = -y_i$ if $0 < \alpha^k_i < C$ or (3b) is satisfied. If $d_i = -y_i$ violates (3b), set $d_i = 0$ and bypass it. From the bottom of the list

*SVM$^{light}$ is available at

http://www-ai.cs.uni-dortmund.de/FORSCHUNG/VERFAHREN/SVM_LIGHT/svm_light.eng.html
sequentially set \( d_i = y_i \) if \( 0 < \alpha_i^k < C \) or (3b) is satisfied. If \( d_i = y_i \) violates (3b), set \( d_i = 0 \) and bypass it. The assignment of \( d_i = -y_i \) and \( y_i \) is done symmetrically until either

(a) \( q/2 \) elements of \( d \) are assigned to be \( -y_i \) from the top and \( q/2 \) elements of \( d \) are assigned to be \( y_i \) from the bottom; or

(b) we cannot find \( d_i = -y_i \) from the top and \( d_i = y_i \) from the bottom at the same time.

3. Elements of \( d \) not considered yet are assigned to be zeros.

Algorithm I.2 will be discussed in more detail later. We mention the working set selection here because it is strongly related to the main topic of this paper: the convergence of the decomposition method.

As the decomposition method finds an optimal solution of a sub-problem (2), the strict decrease of the objective function holds. However, this does not imply that the sequence \( \{\alpha^k\} \) converges to an optimal solution of (1). In fact the convergence issue is not easy and has not been fully understood yet.

The first work on the convergence of the decomposition method is by Chang et al. [3]. They proved the convergence of a more generalized algorithm. However, their working set selection is by a different problem:

\[
\min \nabla f(\alpha^k)^T d \\
0 \leq \alpha_i^k + d_i \leq C, i = 1, \ldots, l, \\
y^T d = 0, \\
|\{d_i \mid d_i \neq 0\}| \leq q.
\]

The main shortcoming is that (4) may not be useful in practice. Unlike Algorithm I.2 for (3), we have not known any comparable method for (4). Note that Algorithm I.2 takes at most \( O(l \ln l) \) or \( O(lq) \) operations that is acceptable for practical implementation.

Then an important progress is by Keerthi and Gilbert [12] where they proved the finite termination of a decomposition method with \( q = 2 \). In [14] the authors showed that the original SMO may not converge so some modifications and improvements were added to SMO. Then [12] intended to prove the finite termination of a generalized SMO algorithm. Incidentally (3) with \( q = 2 \) is a special case of the working set selection proposed in [14] (i.e. modification 2 of SMO in that paper). Thus their proof has covered some existing
practical implementation.

Up to now the only available implementation using $q > 2$ with convergence proofs is discussed in [10]. Instead of using the standard formulation, [10] solves

$$\begin{align*}
\min & \quad \frac{1}{2} \alpha^T (Q + yy^T) \alpha - e^T \alpha \\
0 & \leq \alpha_i \leq C, \quad i = 1, \ldots, l.
\end{align*}$$

(5)

This formulation was proposed and studied by, for example, Friess et al. [9], and Mangasarian and Musicant [16]. (5) is a bound-constrained problem so the working set selection is by the following problem:

$$\begin{align*}
\min & \quad \nabla \bar{f}(\alpha^k)^T d \\
0 & \leq \alpha_i^k + d_i \leq C, i = 1, \ldots, l, \\
|\{d_i \mid d_i \neq 0\}| & \leq q,
\end{align*}$$

(6)

where $\bar{f}(\alpha) \equiv 1/2 \alpha^T (Q + yy^T) \alpha - e^T \alpha$. The convergence follows from the framework in [3]. An important fact is that because of the simpler constraints, (6) can be solved as efficiently as solving (3). To be more precise, the complexity to solve (6) is similar to Algorithm I.2. However, a direct use of (6) did not perform well so [10] finally used a modified way whose convergence is also not clear.

Furthermore, the use of (5) lacks enough theoretical support on generalization properties. We may worry that by removing the linear constraint and adding $1/2(y^T \alpha)^2$ to the objective function, the generalization property is not as good as solving (1). In addition, as more available software follow the implementation of SVM$^{light}$ using (3) (e.g. [6], [21]), the need to prove the convergence with $q > 2$ becomes more emergent. In this paper we will show that Algorithm I.1 using (3) for the working set selection converges.

Next we discuss some possible obstacles while attempting to prove the convergence. In particular, we think the decomposition method of SVM$^{light}$ has two major problems:

1. In each iteration, the decomposition method works only on a subset of variables. Popular optimization methods such as Newton or Quasi Newton consider all variables together in each iteration. In fact if $q$ is small, in each iteration only few coordinates of the variable are updated. Hence the algorithm is like the “coordinate search” or “method of alternating variables” in optimization literature. It has been shown by Powell [20] that such
methods may not always converge. The work in [3] focused on handling this difficulty and a technique to construct a relationship between (4) and the following problem is utilized:

$$\min \ \nabla f (\alpha^k)^T d$$

$$0 \leq \alpha^k_i + d_i \leq C, i = 1, \ldots, l,$$

$$y^T d = 0.$$

(7)

2. **SVM\textsuperscript{light}** uses (3) for the working set selection. Problem (3) follows from the method of feasible directions by Zoutendijk [30]. The original feasible-direction method of Zoutendijk is to consider (3) without restricting the number of nonzero elements:

$$\min \ \nabla f (\alpha^k)^T d$$

$$y^T d = 0, -1 \leq d_i \leq 1, i = 1, \ldots, l,$$

$$d_i \geq 0, \text{ if } \alpha^k_i = 0, \quad d_i \leq 0, \text{ if } \alpha^k_i = C.$$

(8)

The difficulty arises because the convergence of Zoutendijk’s method is not generally guaranteed. The main reason is that $\alpha^k + d$ may not be a feasible point of (1) so the map of search directions is not closed. An example showing that Zoutendijk’s algorithm may not converge is by Wolfe [29] and more discussions are in [1]. This explains why in [3], (6) instead of (3) is considered because (6) guarantees the feasibility of $\alpha^k + d$. To be more precise, the key difficulty is the problem caused by $\alpha_i$ sitting very close to the boundary and having large violation. Another way to see this is that the objective function of (8) is discontinuous at the boundary. This is a rather peculiar situation not associated with traditional optimization algorithms. That is why methods such as Joachims’ decomposition algorithm and SMO by Platt require a very different approach to the proof. The proofs given here and the one in Keerthi and Gilbert use a non-traditional counting argument to prove convergences.

In addition, we note that the original Zoutendijk’s method directly uses $d$ as the search direction for the optimization algorithm. That is, a step size $\lambda$ is decided and $\alpha^k + \lambda d$ becomes the next iterate $\alpha^{k+1}$. This is different from the role of (3) here as $d$ is used only for selecting the working set. Furthermore, in each iteration an exact solution of the sub-problem (2) is obtained. This seems to be a nice property which the original Zoutendijk’s
method lacks of. In [3], such a property was not used as they considered a more general algorithm. For the proof in this paper, we will see that it plays an important role.

The above discussion reveals that the working set selection problem (3) should be deeply investigated. In Section II, we analyze (3) and its solution procedure: Algorithm I.2. Readers who are interested in only the convergence proofs may skip this section. In Section III we sketch the main convergence proof by some figures. Section IV is the main convergence proof. Extensions of the proof to other SVM formulations such as regression and one-class SVM are in Section V. We make conclusions and discussions in Section VI.

II. More Analysis on the Working Set Selection

Though in [11], Joachims has proposed Algorithm I.2 to solve (3), up to now there is no rigorous discussion to justify the use of this algorithm. For example, if the case 2(b) of Algorithm I.2 is encountered first, it is not clear what the practical situation looks like. In this section, we will discuss the details of Algorithm I.2 and demonstrate that it really solves (3). For readers who are interested in only the convergence proofs, you can skip this section and directly go to Section III.

First we give a simple assumption:

Assumption II.1 $C > 0$.

If $C = 0$, the only feasible solution of (1) is $\alpha_i = 0, i = 1, \ldots, l$. In addition, the constraints of (8) implies $d_i = 0$. If Algorithm I.2 is used without Assumption II.1, for $0 = \alpha_i^k = C$ we may end up with $d_i = y_i(-y_i) \neq 0$ which is not a feasible solution of (8). The assumption looks trivial but in our mind we consider the general situation $l_i \leq \alpha_i \leq u_i$, where $l_i$ and $u_i$ are lower and upper bounds, respectively. If $l_i = u_i$, then we can remove the $i$th variable from the original problem easily.

The following theorem shows how Algorithm I.2 solves (8).

Theorem II.2 If the condition 2(a) of Algorithm I.2 is not activated (or $q$ is selected large enough), the algorithm will finally stop at $i_t$ (from the top) and $i_b$ (from the bottom) and one of the following will happen:
1. $y_i \nabla f(\alpha^{k})_{i_t}$ is next to $y_i \nabla f(\alpha^{k})_{i_b}$ in the sorted list.
2. There is one element \( y_i \nabla f(\alpha^k)_i \) between \( y_t \nabla f(\alpha^k)_t \) and \( y_b \nabla f(\alpha^k)_b \) with \( 0 < \alpha_i^k < C \). In addition, when the algorithm stops, \( d \) is an optimal solution of problem (8).

Proof: When Algorithm I.2 stops at \( i_t \), if the next index in the sorted list of \( y_i \nabla f(\alpha^k)_i, i = 1, \ldots, l \) is \( \bar{i}_t \), there are three possible situations:

\[
\begin{align*}
0 < \alpha_i^k &< C, \text{ or} \\
\alpha_i^k &= 0, y_{\bar{i}_t} = -1, \text{ or} \\
\alpha_i^k &= C, y_{\bar{i}_t} = 1.
\end{align*}
\]

Otherwise, we can move down by setting \( d_{\bar{i}_t} = 0 \). Then consider going up from \( i_b \), if the next \( \bar{i}_b \) is not \( \bar{i}_t \) or \( i_t \), it can not satisfy \( 0 < \alpha_i^k < C \), or \( \alpha_i^k = 0, y_{\bar{i}_b} = -1 \), or \( \alpha_i^k = C, y_{\bar{i}_b} = 1 \). Otherwise, (9) implies that \( i_t \) can move down to \( \bar{i}_t \) and then \( i_b \) could move up. Hence \( \bar{i}_b \) must satisfy \( \alpha_i^k = 0, y_{\bar{i}_b} = -1 \) or \( \alpha_i^k = C, y_{\bar{i}_b} = 1 \). However, for this situation, \( i_b \) could move up by setting \( d_{\bar{i}_b} = 0 \). Hence we are sure that there is at most one element between \( i_t \) and \( i_b \). If there is one such element \( \bar{i}_t \) and \( \alpha_i^k = 0, y_{\bar{i}_t} = -1 \), or \( \alpha_i^k = C, y_{\bar{i}_t} = 1 \), \( i_b \) could move up again by assigning \( d_{\bar{i}_t} = 0 \). Therefore, from (9) the only possible situation is to have an element \( i \) between \( i_t \) and \( i_b \) with \( 0 < \alpha_i^k < C \).

Thus we have clarified the situation when the algorithm terminates. Next we will show that when the algorithm stops, the following KKT condition is satisfied so \( d \) is an optimal solution:

\[
\begin{align*}
\nabla f(\alpha^k) &= -by + \lambda_i - \xi_i, \\
y^T d &= 0, \\
\lambda_i(d_i + 1) &= 0, \text{ if } 0 < \alpha_i^k \leq C, \\
\lambda_i d_i &= 0, \text{ if } \alpha_i^k = 0, \\
\xi_i(1 - d_i) &= 0, \text{ if } 0 \leq \alpha_i^k < C, \\
\xi_i d_i &= 0, \text{ if } \alpha_i^k = C, \\
\lambda_i &\geq 0, \xi_i \geq 0, i = 1, \ldots, l.
\end{align*}
\]

If there is an \( \bar{i}_t \) between \( i_t \) and \( i_b \), we select \( b \) such that

\[
y_{\bar{i}_t} \nabla f(\alpha^k)_{\bar{i}_t} + b = 0.
\]
Otherwise, we pick \( b \) such that
\[
y_i \nabla f(\alpha^k)_i + b \geq 0 \quad \text{for the elements before (include) } i_t
\]
\[
y_i \nabla f(\alpha^k)_i + b \leq 0 \quad \text{for the elements after (include) } i_b
\]

Let us consider the case in (11). If \( d_i = -y_i \) and \( y_i = 1 \), by selecting \( \xi_i \equiv 0 \) and \( \lambda_i \equiv \nabla f(\alpha^k)_i \) \( \geq 0 \), (10) is satisfied. The situation is similar for \( y_i = -1 \). If \( d_i = 0 \), there are two possibilities: \( y_i = 1, \alpha^k_i = 0 \) or \( y_i = -1, \alpha^k_i = C \). For the first case, \( \lambda_i \equiv \nabla f(\alpha^k)_i \) \( \geq 0 \), \( \lambda_i d_i \equiv 0 \) and \( \xi_i \equiv 0 \) satisfy (10). The argument for the second case is similar. Furthermore, the same proof can be applied for indices which satisfy \( d_i = y_i \). Thus we have shown that Algorithm I.2 obtains a KKT point. Since (8) is a linear program, a KKT point is an optimal solution.

After the procedure of Algorithm I.2 without activating condition 2(a) for solving (8), we assume that \( i_1, \ldots, i_{mk/2} \) are indices of elements with \( d_i = -y_i \) (in the decreasing order of \( \{y_i \nabla f(\alpha^k)_i\} \)) and \( j_1, \ldots, j_{mk/2} \) are indices of elements with \( d_i = y_i \) (in the increasing order of \( \{y_i \nabla f(\alpha^k)_i\} \)). Then
\[
y_i \nabla f(\alpha^k)_i + b = p_i \geq 0, i = i_1, \ldots, i_{mk/2},
\]
\[
y_i \nabla f(\alpha^k)_i + b = n_i \leq 0, i = j_1, \ldots, j_{mk/2}.
\]
We have
\[
p_{i_1} \geq p_{i_2} \geq \ldots \geq p_{i_{mk/2}} \geq 0 \geq n_{j_{mk/2}} \geq \ldots \geq n_{j_{1}}.
\]
Since \( d_i = -y_i, i = i_1, \ldots, i_{mk/2} \) and \( d_i = y_i, i = j_1, \ldots, j_{mk/2} \),
\[
\nabla f(\alpha^k)_i d_i = (b - p_i), i = i_1, \ldots, i_{mk/2},
\]
\[
\nabla f(\alpha^k)_i d_i = (-b + n_i), i = j_1, \ldots, j_{mk/2}.
\]
Therefore, the optimal objective value of (8) is
\[
\sum_{i = i_1, \ldots, i_{mk/2}} -p_i + \sum_{i = j_1, \ldots, j_{mk/2}} n_i.
\]
Now we are ready to work on problem (3). We will show that by selecting
\[
d_i \equiv \begin{cases} 
-y_i, & i = i_1, \ldots, i_{\min(q, mk)/2}, \\
y_i, & i = j_1, \ldots, j_{\min(q, mk)/2}, \\
0, & \text{otherwise}
\end{cases}
\]
(14)
an optimal solution of (3) is obtained. When \( q \geq m_k \), the solution we just obtained for (8) is a feasible solution of (3). As (3) has a smaller feasible region than (8), its objective value is not smaller. Thus \( d \) defined by (14) is an optimal solution of (3). On the other hand, if \( q < m_k \), we consider the following problem:

\[
\min \quad \nabla f(\alpha^k)^T d \\
y^T d = 0, \quad -1 \leq d_i \leq 1, \quad i = 1, \ldots, l, \\
d_i \geq 0, \quad \text{if } \alpha_i^k = 0, \quad d_i \leq 0, \quad \text{if } \alpha_i^k = C, \\
d_i = 0, \quad \text{if } i \notin \bar{B},
\]

where \( \bar{B} \) is any subset of \( \{1, \ldots, l\} \) containing \( \bar{q} \) elements with \( \bar{q} \leq q \). Now \( \bar{B} \) is fixed so (15) is reduced to a form of (8) whose number of variables is \( \bar{q} \). Hence the same procedure of Algorithm I.2 without Step 2(a) could be applied to solve (15). If an optimal solution is \( d_i = -y_i, i = r_1, \ldots, r_{\bar{q}/2}, \quad d_i = y_i, i = s_1, \ldots, s_{\bar{q}/2} \) (\( \bar{q} \leq \bar{q} \)), and \( d_i = 0 \) otherwise, then the optimal objective value of (15) is

\[
\left( -\sum_{i=r_1,\ldots,\bar{q}/2} y_i \nabla f(\alpha^k)_i + \sum_{i=s_1,\ldots,\bar{q}/2} y_i \nabla f(\alpha^k)_i \right),
\]

which is greater or equal to

\[
\left( -\sum_{i=1,\ldots,\bar{q}/2} y_i \nabla f(\alpha^k)_i + \sum_{i=1,\ldots,\bar{q}/2} y_i \nabla f(\alpha^k)_i \right),
\]

as we sort \( y_i \nabla f(\alpha^k)_i \) in a decreasing order. Since \( \hat{q} \leq \bar{q} \leq q \), \( d \) defined in (14) is an optimal solution of (3). The following theorem concludes the validity of Algorithm I.2 for (3):

**Theorem II.3** If \( q \) is an even positive integer, Algorithm I.2 returns an optimal solution of (3) and

\[
\frac{l}{q} \text{(optimal objective value of (3))} \leq \text{optimal objective value of (8)}. \quad (16)
\]

We then show the relation between the working set selection problem (3) and the original optimization problem (1):

**Theorem II.4** The optimal objective value of (3) is zero if and only if \( \alpha \) is an optimal solution of (1).
Proof: A basic property of Zoutendijk’s method is that the optimal objective value of (8) is zero if and only if $\alpha$ is an optimal solution of (1) (see, for example, [1]). Since (3) has a smaller feasible region than (8), if the optimal objective value of (3) is zero, the optimal solution of (8) is also zero. Therefore, $\alpha$ is an optimal solution of (1).

On the other hand, if $\alpha$ is an optimum of (1), with Lemma II.3, the optimal objective value of (3) is zero.

There are different methods for the analysis in this section. For example, in [4], the authors modified (3) to $\nu$-SVM problems [25] where they used a recursive approach to show the validity of Algorithm I.2.

III. OUTLINE OF THE CONVERGENCE PROOF

The convergence of Algorithm I.1 using problem (3) is the main result of this paper. As the proof involves with several complicated lemmas and theorems, in this section we give an outline of the proof. Using some informal terms and figures, we explain some key ideas behind the proof.

First we discuss some observations which help to prove the convergence. If $\hat{\alpha}$ is an optimal solution of (1), it satisfies the following KKT condition: there is a number $b$ such that

$$\nabla f(\hat{\alpha})_i + by_i \geq 0 \quad \text{if } \hat{\alpha}_i = 0,$$
$$\nabla f(\hat{\alpha})_i + by_i \leq 0 \quad \text{if } \hat{\alpha}_i = C,$$
$$\nabla f(\hat{\alpha})_i + by_i = 0 \quad \text{if } 0 < \hat{\alpha}_i < C. \quad (17)$$

For any scalar $\alpha_i$, we can consider two situations

$$0 < \alpha_i < C \quad \text{or} \quad (\alpha_i = C \text{ and } y_i = 1) \quad \text{or} \quad (\alpha_i = 0 \text{ and } y_i = -1), \quad (18)$$
$$0 < \alpha_i < C \quad \text{or} \quad (\alpha_i = C \text{ and } y_i = -1) \quad \text{or} \quad (\alpha_i = 0 \text{ and } y_i = 1). \quad (19)$$

Then the KKT condition (17) can be rewritten as

$$y_i \nabla f(\hat{\alpha})_i + b \geq 0 \quad \text{if } \hat{\alpha}_i \text{ satisfies (19)},$$
$$y_i \nabla f(\hat{\alpha})_i + b \leq 0 \quad \text{if } \hat{\alpha}_i \text{ satisfies (18)}. \quad (20)$$
Note that (18) ((19)) is the condition in Algorithm I.2 where $\alpha_i^k$ can be a candidate for selection from the top (bottom) of the sorted list of $y_i \nabla f(\alpha_i^k), i = 1, \ldots, l$. In the following we shall refer a variable $\alpha_i$ as a

“top” candidate: if it satisfies (18),

“top only” candidate: if it satisfies ($\alpha_i = C$ and $y_i = 1$) or ($\alpha_i = 0$ and $y_i = -1$),

“bottom” candidate: if it satisfies (19), or

“bottom only” candidate: if it satisfies ($\alpha_i = C$ and $y_i = -1$) or ($\alpha_i = 0$ and $y_i = 1$).

From Assumption II.1, $C > 0$ so the following two statements are equivalent:

$\alpha_i$ is a “top only” candidate $\equiv \alpha_i$ is not a “bottom” candidate.

Therefore, once $\alpha_i^k$ is a “top only” candidate, next time when it is selected, in Algorithm I.2, it must be picked from the top of the sorted list.

\[
\begin{array}{cccccc}
\alpha^k & \alpha^{k+1} & \bar{\alpha} & \alpha^k & \alpha^{k+1} & \bar{\alpha} \\
i_1(\text{top}) & i_1(\text{top}) & i_1(\text{top}) & i_1 \in B & i_1(\text{top}) & i_1 \\
i_2(\text{bottom}) & i_2(\text{bottom}) & i_2(\text{bottom}) & i_2 \in B & i_2(\text{bottom}) & i_2 \\
\cdots & \rightarrow & \rightarrow & \end{array}
\]

(a) A convergent subsequence (Lemmas IV.3 and IV.4(a))
(b) An impossible situation (Lemma IV.4(b))

Fig. 1. $y_i \nabla f(\alpha_i)$ in the order of the sorted list of $y_i \nabla f(\bar{\alpha}_i), i = 1, \ldots, l$

It can be clearly seen that the KKT condition (20) implies that all “top” candidates have the same or smaller $y_i \nabla f(\hat{\alpha}_i)$ than “bottom” candidates. Therefore, when applying Algorithm I.2 to problem (3) of an optimal solution, except those elements with $y_i \nabla f(\hat{\alpha}_i) + b = 0$, we cannot do any selection.

For free variables, their $y_i \nabla f(\hat{\alpha}_i)$ are equal. On the other hand, if $\hat{\alpha}_i$ are at bounds, their $y_i \nabla f(\hat{\alpha}_i)$ are usually different. This leads us to suspect that in final iterations, all bounded $\alpha_i$’s associated $y_i \nabla f(\alpha_i)$ are already in correct places of the sorted list of $y_i \nabla f(\bar{\alpha}_i), i = 1, \ldots, l$.

Of course it is possible that $\nabla f(\hat{\alpha}_i) + b y_i = 0$ even if $\hat{\alpha}_i$ is at a bound. This is the so called “degenerate” case in optimization terminology. For degenerate or free variables,
\[ \alpha^k \quad \alpha^k \quad \alpha^{k+1} \quad \alpha^{k+2(l+1)} \quad \bar{\alpha} \]

<table>
<thead>
<tr>
<th>( I_1 )</th>
<th>( I_1 ) (bottom only)</th>
<th>( \cdots )</th>
<th>( I_1 ) (top)</th>
<th>( \cdots )</th>
</tr>
</thead>
<tbody>
<tr>
<td>( i_1 ) (top)</td>
<td>( i_1 ) (top) ( \in B )</td>
<td>( \cdots )</td>
<td>( i_1 ) (top)</td>
<td>( \cdots )</td>
</tr>
<tr>
<td>( \cdots )</td>
<td>( \cdots )</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>( i_2 ) (bottom)</td>
<td>( i_2 ) (bottom) ( \in B )</td>
<td>( i_2 ) (bottom)</td>
<td>( \cdots )</td>
<td></td>
</tr>
<tr>
<td>( I_2 )</td>
<td>( I_2 ) (top only)</td>
<td>( \cdots )</td>
<td>( I_2 )</td>
<td></td>
</tr>
</tbody>
</table>

Fig. 2. A counting process on \( I_1 \) and \( I_2 \) (Theorem IV.5)

\( y_i \nabla f(\hat{\alpha}) \) are all equal. We will focus on analyzing this group of variables. Indeed we will show that for any limit point of a convergent subsequence, only indices from this particular group are still under consideration.

Next we outline the proof. We consider any convergent subsequence \( \{\alpha^k\}, k \in \mathcal{K} \) and \( \bar{\alpha} \equiv \lim_{k \to \infty, k \in \mathcal{K}} \alpha^k \). We prove that for any given positive integer \( s \), the sequence \( \{\alpha^{k+s}\}, k \in \mathcal{K} \) converges to \( \bar{\alpha} \). Therefore, if \( \bar{\alpha}_i \) is a “top” (“bottom”) candidate, then after \( k \in \mathcal{K} \) is large enough, \( \alpha_i^k, \alpha_i^{k+1}, \cdots, \alpha_i^{k+s} \) are all “top” (“bottom”) candidates. These results will be proved in Lemmas IV.3 and IV.4(a). An illustration is in Figure 1(a), where the vertical line on the right represents the sorted list \( y_i \nabla f(\bar{\alpha})_i, i = 1, \ldots, l \) (in the decreasing order) and lines on the left are corresponding values of \( y_i \nabla f(\alpha^k)_i \) to \( y_i \nabla f(\alpha^{k+s})_i \).

We then prove that a situation like Figure 1(b) cannot happen. That is, if

\[ y_{i_1} \nabla f(\bar{\alpha})_{i_1} > y_{i_2} \nabla f(\bar{\alpha})_{i_2}, \]

then after \( k \in \mathcal{K} \) is large enough, for any \( \bar{k} \in \{k, k + 1, \ldots, k + s - 1\} \), if \( i_1 \) and \( i_2 \) are both in the working set of the \( \bar{k} \)th iteration, it is impossible to have \( \alpha_i^{k+1} \) and \( \alpha_i^{k+1} \) are “top” and “bottom” candidates at the same time.

The final part of the proof comes from Figure 2 (Theorem IV.5). Assume \( i_1(i_2) \) is the first “top” (“bottom”) candidate of the sorted list of \( y_i \nabla f(\bar{\alpha})_i, i = 1, \ldots, l \) and \( I_1(I_2) \) is the set not lower (higher) than \( i_1(i_2) \). If

\[ y_{i_1} \nabla f(\bar{\alpha})_{i_1} > y_{i_2} \nabla f(\bar{\alpha})_{i_2}, \]

we prove that after \( k \in \mathcal{K} \) is large enough, there is a \( \bar{k} \in \{k, k + 1, \ldots, k + 2l\} \) such that \( \alpha_{i_1}^{k} \) (\( \alpha_{i_2}^{k} \)) has only “bottom only” (“top only”) elements. Therefore, at the \( \bar{k} \)th iteration, both
$i_1$ and $i_2$ must be selected as they are the first “top” and “bottom” elements, respectively (from Figure 1(a)). However, at the $(\bar{k} + 1)\text{st}$ iteration, from Figure 1(a), we know that $i_1$ and $i_2$ are “top” and “bottom” elements again. This violates the results in Figure 1(b). Therefore, the assumption (21) is wrong so at $\bar{\alpha}$, elements selected by (3) have the same $y_i \nabla f(\bar{\alpha})$. With this result we can show that the working set selection problem at $\bar{\alpha}$ has zero optimal objective value. Then from Theorem II.4, $\bar{\alpha}$ is an optimal solution of (1).

Thus the main effort of Theorem IV.5 is on a counting process to show that in at most $2l$ iterations, all elements in $I_1$ and $I_2$ become “bottom only” and “top only,” respectively.

IV. Convergence Proofs

In this section we prove the convergence of Algorithm I.1 using problem (3) for the working set selection (i.e. the algorithm used by SVM$^{light}$). If Algorithm I.1 stops in finite number of iterations, from Step 2, $\alpha^k$ is already an optimum. Hence here we consider the case where Algorithm I.1 takes infinite iterations. First we make an assumption:

**Assumption IV.1** The matrix $Q$ satisfies

$$\min_I(\min(eig(Q_{II}))) > 0,$$

where $I$ is any subset of $\{1, \ldots, l\}$ with $|I| \leq q$, $Q_{II}$ is a square sub-matrix of $Q$, and $\min(eig(\cdot))$ is the smallest eigenvalue of a matrix.

If $Q$ is positive definite, then Assumption IV.1 is true. For example, if the RBF kernel $K(x_i, x_j) = e^{\|x_i - x_j\|^2}$ is used and all $x_i \neq x_j$, from [17], $Q$ is positive definite. Since practically $q$ is selected as a small number ($\leq 100$), if data are mapped into higher dimensional spaces, $Q$ tends to be positive definite so in general Assumption IV.1 holds.

The following lemma shows the sufficient decrease of $f(\alpha)$:

**Lemma IV.2**

$$f(\alpha^{k+1}) \leq f(\alpha^k) - \frac{\sigma}{2} \|\alpha^{k+1} - \alpha^k\|^2,$$  \hspace{1cm} (22)

where $\sigma = \min_I(\min(eig(Q_{II})))$. 

**Proof:** Assume $B$ is the working set at the $k$th iteration and $N \equiv \{1, \ldots, l\} \setminus B$. If we define $s \equiv \alpha^{k+1} - \alpha^k$, then $s_N = 0$ and

$$
\begin{align*}
& f(\alpha^{k+1}) - f(\alpha^k) \\
& = \frac{1}{2} s^T Q s + s^T Q \alpha^k - e^T s \\
& = \frac{1}{2} s_B^T Q_B s_B + s_B^T (Q \alpha^k)_B - e_B^T s_B.
\end{align*}
$$

That is, in the $k$th iteration, we solve the following problem with the variable $s_B$:

$$
\begin{align*}
\min & \quad \frac{1}{2} s_B^T Q_B s_B + s_B^T (Q \alpha^k)_B - e_B^T s_B \\
\text{s.t.} & \quad 0 \leq (\alpha^k + s)_i \leq C, i \in B, \\
& \quad y_B^T s_B = 0,
\end{align*}
$$

which is a different representation of (2). The KKT condition of (24) shows that there is a $b^{k+1}$ such that

$$
\begin{align*}
(Q(\alpha^k + s))_i - 1 + b^{k+1} y_i &= 0 & \text{if } 0 < \alpha^k_i + s_i < C, i \in B, \\
(Q(\alpha^k + s))_i - 1 + b^{k+1} y_i &\geq 0 & \text{if } \alpha^k_i + s_i = 0, i \in B, \\
(Q(\alpha^k + s))_i - 1 + b^{k+1} y_i &\leq 0 & \text{if } \alpha^k_i + s_i = C, i \in B.
\end{align*}
$$

Define $F \equiv \{i \mid 0 < \alpha^k_i + s_i < C, i \in B\}$ and $A \equiv \{i \mid \alpha^k_i + s_i = 0 \text{ or } C, i \in B\}$. We have $B = F \cup A$ and from (27),

$$
(Q \alpha^k)_F = -(Q s)_F + e_F - b^{k+1} y_F = -Q_F s_F - Q_F s_A + e_F - b^{k+1} y_F
$$

With (28), the last two terms of (23) become

$$
\begin{align*}
& s_B^T (Q \alpha^k)_B - e_B^T s_B \\
& = s_F^T (Q \alpha^k)_F - e_F^T s_F + s_A^T ((Q(\alpha^k + s))_A + b^{k+1} y_A - e_A) - s_A^T (Q s)_A - b^{k+1} y_A^T s_A \\
& = s_F^T (Q \alpha^k)_F - e_F^T s_F + s_A^T ((Q(\alpha^k + s))_A + b^{k+1} y_A - e_A) - s_A^T (Q A s_F + Q A s_A) - b^{k+1} y_A^T s_A \\
& = -s_F^T Q_F s_F - s_A^T ((Q(\alpha^k + s))_A + b^{k+1} y_A - e_A) - s_A^T (Q A s_F + Q A s_A).
\end{align*}
$$
If \( \alpha_i^k + s_i = 0 \), then \( s_i \leq 0 \) and if \( \alpha_i^k + s_i = C \), then \( s_i \geq 0 \). Hence from (26) and (27)

\[
s_A^T((Q(\alpha^k + s))_A + b^{k+1} y_A - e_A) \leq 0.
\]

With (29), (30),

\[
\frac{1}{2} s_B^T Q_{BB} s_B = \frac{1}{2} s_F^T Q_{FF} s_F + s_F^T Q_{FA} s_A + \frac{1}{2} s_A^T Q_{AA} s_A,
\]

and \( y_B^T s_B = 0 \), (23) becomes

\[
-\frac{1}{2} s_F^T Q_{FF} s_F - \frac{1}{2} s_A^T Q_{AA} s_A - s_F^T Q_{FA} s_A + s_A^T((Q(\alpha^k + s))_A + b^{k+1} y_A - e_A)
\]

\[
\leq - \frac{1}{2} \begin{bmatrix} s_F^T & s_A^T \end{bmatrix} \begin{bmatrix} Q_{FF} & Q_{FA} \\ Q_{AF} & Q_{AA} \end{bmatrix} \begin{bmatrix} s_F \\ s_A \end{bmatrix}
\]

\[
\leq - \frac{\sigma}{2} \|s_B\|^2 = - \frac{\sigma}{2} \|s\|^2.
\]

From now on we consider any convergent subsequence \( \{\alpha^k\}, k \in \mathcal{K} \) and \( \lim_{k \to \infty, k \in \mathcal{K}} \alpha^k = \bar{\alpha} \). We then have the following lemma:

**Lemma IV.3** For any given positive integer \( s \), the sequence \( \{\alpha^{k+s}\}, k \in \mathcal{K} \) converges to \( \bar{\alpha} \). In addition, \( \{y_i \nabla f(\alpha^{k+s})_i\} \) converges to \( y_i \nabla f(\bar{\alpha})_i \), for \( i = 1, \ldots, l \).

*Proof:* First we know that \( \{f(\alpha^k)\} \) is a decreasing sequence. Since \( 0 \leq \alpha_i \leq C \), \( i = 1, \ldots, l \), the feasible region of (1) is a compact set. Thus we know that \( \{f(\alpha^k)\} \) converges to a finite number.

Then for the subsequence \( \{\alpha^{k+1}\}, k \in \mathcal{K} \), from Lemma IV.2 we have

\[
\lim_{k \to \infty} \|\alpha^{k+1} - \bar{\alpha}\|
\]

\[
\leq \lim_{k \to \infty} (\|\alpha^{k+1} - \alpha^k\| + \|\alpha^k - \bar{\alpha}\|)
\]

\[
\leq \lim_{k \to \infty} \left( \frac{2}{\sigma} (f(\alpha^k) - f(\alpha^{k+1})) + \|\alpha^k - \bar{\alpha}\| \right)
\]

\[
= 0.
\]

Thus

\[
\lim_{k \to \infty, k \in \mathcal{K}} \alpha^{k+1} = \bar{\alpha}.
\]
From \( \{\alpha^k\} \) we can prove \( \lim_{k \to \infty, k \in K} \alpha^{k+2} = \bar{\alpha} \) too. Therefore, \( \lim_{k \to \infty, k \in K} \alpha^{k+s} = \bar{\alpha} \) for any given \( s \).

The results on \( \{y_i \nabla f(\alpha^{k+s})_i\} \) follows from the continuity of \( \nabla f(\alpha) \).

We then need a technical lemma:

**Lemma IV.4** Let \( \bar{\alpha} \) be as in Lemma IV.3.

(a) If \( \bar{\alpha}_i \) satisfies (18) ((19)), then for any given positive integer \( s \), after \( k \in K \) is large enough, \( \alpha^k_i, \alpha^{k+1}_i, \ldots, \alpha^{k+s}_i \) all satisfy (18) ((19)). In other words, if \( \bar{\alpha}_i \) is a “top” (“bottom”) candidate, then after \( k \in K \) is large enough, \( \alpha^k_i, \alpha^{k+1}_i, \ldots, \alpha^{k+s}_i \) are all “top” (“bottom”) candidates.

(b) In addition, if

\[
y_i \nabla f(\bar{\alpha})_i > y_{i_2} \nabla f(\bar{\alpha})_{i_2},
\]

then after \( k \in K \) is large enough, for any \( \bar{k} \in \{k, k+1, \ldots, k+s-1\} \), if \( i_1 \) and \( i_2 \) are both in the working set of the \( \bar{k} \)th iteration, it is impossible to have \( \alpha^{\bar{k}+1}_{i_1} \) and \( \alpha^{\bar{k}+1}_{i_2} \) satisfy (18) and (19), respectively. In other words, \( \alpha^{\bar{k}+1}_{i_1} \) and \( \alpha^{\bar{k}+1}_{i_2} \) cannot be top and bottom candidates at the same time.

**Proof:** The first result immediately follows from Assumption II.1, Lemma IV.3, and the definition of (18) and (19).

For the second result of this lemma, we assume that it is possible that both \( \alpha^{\bar{k}+1}_{i_1} \) and \( \alpha^{\bar{k}+1}_{i_2} \) satisfy (18) and (19), respectively. Since \( \alpha^{\bar{k}+1}_B \) is an optimal solution of (2), from the KKT condition of the sub-problem (2) and a similar form of (20), if \( \alpha^{\bar{k}+1}_{i_1} \) satisfies (18), there is a \( b^{\bar{k}+1} \) such that

\[
y_{i_1} \nabla f(\alpha^{\bar{k}+1})_{i_1} + b^{\bar{k}+1} \leq 0.
\]

On the other hand, if \( \alpha^{\bar{k}+1}_{i_2} \) satisfies (19), then

\[
y_{i_2} \nabla f(\alpha^{\bar{k}+1})_{i_2} + b^{\bar{k}+1} \geq 0.
\]

Thus (33) and (34) imply

\[
y_{i_1} \nabla f(\alpha^{\bar{k}+1})_{i_1} \leq y_{i_2} \nabla f(\alpha^{\bar{k}+1})_{i_2}
\]

which contradicts to (32) when \( \bar{k} \) is large enough.
Finally, the main theorem is as follows:

**Theorem IV.5** Any limit point of \( \{\alpha^k\} \) is a global minimum of (1).

**Proof:** Assume \( \bar{\alpha} \) is the limit point of any convergent subsequence \( \{\alpha^k\}, k \in K \). If \( \bar{\alpha} \) is not an optimal solution of (1), from Theorem II.4, the following problem has a nonzero solution:

\[
\begin{align*}
\min & \quad \nabla f(\bar{\alpha})^T \mathbf{d} \\
-1 \leq d_i & \leq 1, y^T \mathbf{d} = 0, \\
 d_i & \geq 0, \text{ if } \bar{\alpha}_i = 0, \quad d_i \leq 0, \text{ if } \bar{\alpha}_i = C, \\
\left| \{d_i \mid d_i \neq 0\} \right| & \leq q.
\end{align*}
\]

(35)

If we can prove that only elements with the same \( y_i \nabla f(\bar{\alpha})_i \) can have nonzero \( d_i \), by assuming \( B \) contains such indices, then

\[
\nabla f(\bar{\alpha})^T \mathbf{d} = \sum_{i \in B} y_i^2 \nabla f(\bar{\alpha})_i d_i = (y_i \nabla f(\bar{\alpha})_i) y_B^T \mathbf{d}_B = 0.
\]

This contradicts to the assumption that (35) has a nonzero solution. Hence \( \bar{\alpha} \) is an optimal solution.

Therefore, in the rest of this proof we will show that when solving (35), only elements with the same \( y_i \nabla f(\bar{\alpha})_i \) can have nonzero \( d_i \). Assume \( i_1 \) (\( i_2 \)) is the first element selected from the top (bottom) of the sorted list of \( y_i \nabla f(\bar{\alpha})_i, i = 1, \ldots, l \). We claim that

\[
y_{i_1} \nabla f(\bar{\alpha})_i = y_{i_2} \nabla f(\bar{\alpha})_i.
\]

If the result is wrong, then

\[
y_{i_1} \nabla f(\bar{\alpha})_i > y_{i_2} \nabla f(\bar{\alpha})_i.
\]

(36)

Define

\[
I_1 \equiv \{i \mid y_i \nabla f(\bar{\alpha})_i \geq y_{i_1} \nabla f(\bar{\alpha})_{i_1}\}
\]

and

\[
I_2 \equiv \{i \mid y_i \nabla f(\bar{\alpha})_i \leq y_{i_2} \nabla f(\bar{\alpha})_{i_2}\}.
\]
From (36), $I_1 \cap I_2 = \emptyset$.

Since $i_1$ ($i_2$) is the first element selected from the top (bottom) of the sorted list, $\bar{\alpha}_{i_1}$ ($\bar{\alpha}_{i_2}$) satisfies (18) ((19)). After $k \in \mathcal{K}$ is large enough, from Lemma IV.4, $\alpha^k_{i_1}, \alpha^{k+1}_{i_1}, \ldots, \alpha^{k+2l}_{i_1}$ are all “top” candidates, where $l$ is the length of each vector $\alpha$ (i.e. the number of variables of (1)). In addition, $\alpha^k_{i_2}, \alpha^{k+1}_{i_2}, \ldots, \alpha^{k+2l}_{i_2}$ are all “bottom” candidates. Then in each $\bar{k}$ of $k$th, $(k+1)$st, $\ldots$, $(k+2l-1)$st iterations, $i_1$ and $i_2$ can not both be selected because of (36) and Lemma IV.4.

We then claim that if $i_1$ is not selected at the $\bar{k}$th iteration, then all “top” candidates selected are from $I_1$. Since $\bar{k}$ is large enough, for any $\alpha^k_{i_1}, i \notin I_1$, which is a “top” candidate,

$$y_i \nabla f(\bar{\alpha})_i < y_{i_1} \nabla f(\bar{\alpha})_{i_1}$$

implies that $i$ can not be chosen earlier then $i_1$. Similarly, if $i_2$ is not selected at the $\bar{k}$th iteration, then all “bottom” candidates selected are from $I_2$.

Now for the $\bar{k}$th iteration, we consider three situations:

Case 1: Neither $i_1$ nor $i_2$ is selected: Then all “top” (“bottom”) candidates selected are in $I_1(I_2)$. For any $i \in I_1$ and $j \in I_2$ selected in the $\bar{k}$th iteration, from Lemma IV.4 (b), at the next iteration, either $\alpha^k_{i_1}$ becomes a “bottom only” element or $\alpha^k_{i_2}$ becomes a “top only” element. Therefore, there are two cases to consider:

Case 1-1: All elements selected from $I_1$ become “bottom only”: Then the number of “bottom only” variables in $I_1$ is increased by at least one. On the other hand, since $I_1 \cap I_2 = \emptyset$ and from the assumption of case 1, all variables selected from $I_2$ are “bottom” elements. Hence the number of “top only” variables in $I_2$ is at least the same.

Case 1-2: All elements selected from $I_2$ become “top only”: Similarly, the number of “top only” variables in $I_2$ is increased by at least one, while the number of “bottom only” variables in $I_1$ is at least the same.

Case 2: Only $i_1$ is selected: As $i_2$ is not selected, all “bottom” elements selected are in $I_2$. Since $i_1$ is selected and $\alpha^k_{i_1}$ is a “top” candidate, all “bottom” elements selected in $I_2$ become “top only.” Therefore, the number of “top only” variables in $I_2$ increases by at least one. On the other hand, the number of “bottom only” variables in $I_1$ is at least the same.
Case 3: Only $i_2$ is selected: Similar to case 2, the number of “bottom only” variables in $I_1$ increases at least one and the number of “top only” variables in $I_2$ is at least the same.

Therefore, in at most $l$ iterations, either all elements in $I_1$ become “bottom only” or all elements in $I_2$ become “top only.” If $I_1$ reaches “bottom only” first, from Assumption II.1, for later iterations, elements in $I_1$ are not “top” candidates so $i_1$ must be selected. Therefore, we only have case 2 left. Then after at most another $l$ iterations, all $I_2$ are “top only.” Therefore, we must have a $\bar{k} \in \{k, k + 1, \ldots, k + 2l\}$ such that both $i_1$ and $i_2$ are selected. This contradicts to Lemma IV.4. Hence the proof is complete.

Under some conditions (for example, $Q$ is positive definite), (1) has a unique solution. Hence $\{\alpha^k\}$ is a globally convergent sequence whose limit point is this unique solution. In Burges and Crisp [2], there are discussions on conditions under which the SVM solution is unique.

V. Extensions

Consider a general problem with the following form:

\[
\min \frac{1}{2} \alpha^T Q \alpha + p^T \alpha \\
y^T \alpha = \Delta, \\
l_i \leq \alpha_i \leq u_i, i = 1, \ldots, l,
\]

where $-\infty < l_i < u_i < \infty, i = 1, \ldots, l$, $Q$ is any symmetric positive semi-definite matrix satisfying Assumption IV.1, and $y_i = \pm 1, i = 1, \ldots, l$. The convergence proof described in the previous section is still valid if Algorithm I.1 with the following generalized working set selection is used for solving (37):

\[
\min \nabla f(\alpha^k)^T d \\
y^T d = 0, \quad -1 \leq d_i \leq 1, i = 1, \ldots, l, \\
d_i \geq 0, \text{ if } (\alpha^k)_i = l_i, \quad d_i \leq 0, \text{ if } (\alpha^k)_i = u_i, \\
\{|d_i \mid d_i \neq 0\} \leq q.
\]

It can be seen that $y_i = \pm 1$ plays an important role here. Algorithm I.2 is not valid for solving (38) if this condition does not hold. In addition, in the convergence proof we
specifically utilize many properties of Algorithm I.2 (e.g. we consider the sorted list of 
\(y_i \nabla f(\alpha)_i\)) so the condition \(y_i = \pm 1\) is also used. In [12], the authors handled a more
generalized problem where the only restriction on \(y_i\) is \(y_i \neq 0\).

Problem (37) covers most SVM formulations. For example, given a set of data points
\[\{(x_1, z_1), \ldots, (x_l, z_l)\}\] such that \(x_i \in \mathbb{R}^n\) is an input and \(z_i \in \mathbb{R}^1\) is a target output, the
usual form of support vector regression is as follows:

\[
\min \frac{1}{2}(\alpha - \alpha^*)^T Q(\alpha - \alpha^*) + \epsilon \sum_{i=1}^{l} (\alpha_i + \alpha_i^*) + \sum_{i=1}^{l} z_i(\alpha_i - \alpha_i^*)
\]

\[
\sum_{i=1}^{l} (\alpha_i - \alpha_i^*) = 0, 0 \leq \alpha_i, \alpha_i^* \leq C, i = 1, \ldots, l,
\]

(39)

where \(Q_{ij} = \phi(x_i)^T \phi(x_j)\).

We can rewrite (39) as

\[
\min \frac{1}{2} \begin{bmatrix} \alpha^T, (\alpha^*)^T \end{bmatrix} \begin{bmatrix} Q & -Q \\ -Q & Q \end{bmatrix} \begin{bmatrix} \alpha \\ \alpha^* \end{bmatrix} + \begin{bmatrix} \epsilon e^T + z^T, \epsilon e^T - z^T \end{bmatrix} \begin{bmatrix} \alpha \\ \alpha^* \end{bmatrix}
\]

\[
y^T \begin{bmatrix} \alpha \\ \alpha^* \end{bmatrix} = 0, 0 \leq \alpha_i, \alpha_i^* \leq C, i = 1, \ldots, l,
\]

(40)

where \(y\) is a \(2l\) by 1 vector with \(y_i = 1, i = 1, \ldots, l\) and \(y_i = -1, i = l + 1, \ldots, 2l\). (40) is
in the form of (37) so Algorithms I.1 and I.2 can be applied.

However, using Algorithms I.1 and I.2 for (40) is a little different from existing decom-
position methods for regression. Note that though (39) is a problem with \(2l\) variables, it
has very special structures. For example, the KKT condition implies that at an optimal
solution of (39), \(\alpha_i \alpha_i^* = 0\). Early work on SVM regression (e.g. [26], [13], [15], [8]) all
tried to take advantage of these structures and focused on problem (39). Except [15]
they mainly consider selecting two elements as the working set in each iteration. Some
characteristics of their methods are:

1. In each iteration, two indices \(i_1\) and \(i_2\) are selected from \(\{1, \ldots, l\}\).
2. To keep \(\alpha_i^k(\alpha^*)^k = 0\), they solve a sub-problem with four variables \(\alpha_{i_1}, \alpha_{i_2}, \alpha_{i_1}^*,\) and \(\alpha_{i_2}^*\).

That is, it is like that they use \(q = 4\) in Algorithm I.1 with a different working set selection
from Algorithm I.2.
If Algorithms I.1 and I.2 with \( q = 2 \) are directly used for (40), two indices are selected from \( \{1, \ldots, 2l\} \) and a sub-problem (2) with two variables is solved. The advantage of working on (40) is that a generalized implementation can be directly used for both classification and regression. However, a possible shortcoming is that special structures of (39) are not considered so there may have computational overheads. Surprisingly we will show that \( \alpha_i^k(\alpha^*)^k_i = 0 \) still holds if Algorithms I.1 and I.2 are directly applied to solve (40).

Another issue is on the Hessian \( \bar{Q} = [Q - Q Q] \) of the objective function of (40). Now \( \bar{Q} \) is only positive semidefinite so it is unlikely that Assumption IV.1 can be true. In the following theorem we will show that if only \( Q \) instead of \( \bar{Q} \) satisfies Assumption IV.1, the convergence for solving (40) follows.

**Theorem V.1** If Algorithms I.1 and I.2 are used for solving (40) and the initial solution is zero, then \( \alpha_i^k(\alpha^*)^k_i = 0, i = 1, \ldots, l \) for all \( k \). In addition, if \( Q \) satisfies Assumption IV.1, \( \left[ \frac{\alpha_i^k}{(\alpha^*)_k} \right] \) converges to an optimal solution of (39).

**Proof:**

We prove the first result by the mathematical induction. It is true that if the initial solution is zero, for the first iteration, \( \alpha_1^1(\alpha^*)_1^1 = 0, i = 1, \ldots, l \). Assume the result is true for the \( k \)th iteration and we will prove \( \alpha_i^{k+1}(\alpha^*)^{k+1}_i = 0, i = 1, \ldots, l \).

We consider three situations in the \( k \)th iteration:

1. In Algorithm I.2, both \( i \) and \( i + l \) are selected in the working set: Then from the KKT condition of the sub-problem (2), \( \alpha_i^{k+1}(\alpha^*)_i^{k+1} = 0 \).

2. Only \( i \) but not \( i + l \) is selected in the working set and \( \alpha_i^k = 0 \): For this case, \( d_i = 1 \) after solving (3). Since \( y_id_i = 1 \), we realize that in Algorithm I.2, the index \( i \) is selected from the bottom of the sorted list of \( y_i \nabla f(\alpha^k, (\alpha^*)^k)_i, i = 1, \ldots, 2l \). However, we also have

\[
y_i \nabla f(\alpha^k, (\alpha^*)^k)_i = (Q(\alpha^k - (\alpha^*)^k))_i + \epsilon + z_i \geq (Q(\alpha^k - (\alpha^*)^k))_i - \epsilon + z_i = y_{i+l} \nabla f(\alpha^k, (\alpha^*)^k)_{i+l}.
\]

In other words, index \( i + l \) is closer than \( i \) to the bottom of the sorted list. Therefore, if \( i + l \) is not selected, index \( i + l \) is not a “bottom” candidate so \( (\alpha^*)^k_i \) does not satisfy (19).

As \( y_{i+l} = -1 \), \( (\alpha^*)_i = 0 \). Since \( i + l \) is not selected in the \( k \)th iteration, \( (\alpha^*)^{k+1}_i = 0 \) so \( \alpha_i^{k+1}(\alpha^*)^{k+1}_i = 0 \).
3. Only $i$ but not $i+l$ is selected in the working set and $\alpha^k_i > 0$: Then $\alpha^k_i (\alpha^*)^k_i = 0$ implies $(\alpha^*)^k_i = 0$. Therefore, $(\alpha^*)^{k+1} = 0$ and $\alpha^{k+1}_i (\alpha^*)^{k+1}_i = 0$.

When only $i+l$ but not $i$ is selected, the situation is similar. Thus we have finished the proof that $\alpha^k_i (\alpha^*)^k_i = 0$, $i = 1, \ldots, l$, for all $k$.

Next we switch to the second goal of this theorem. Now the Hessian of the objective function of (40) is $\bar{Q} = \begin{bmatrix} Q & -Q \\ -Q & Q \end{bmatrix}$. We remember that Assumption IV.1 is needed near Eq. (31) in the proof of Lemma IV.2. If we can prove
\[
-\frac{1}{2} s^T_B \bar{Q} B B^T s_B \leq -\frac{\sigma}{2} \|s_B\|^2,
\]
where $\sigma$ is related only to $Q$, then a condition on $Q$ instead of $\bar{Q}$ is sufficient for the convergence. Thus the main task is to prove that (41) is true.

We define the following disjoint index sets:

\[
B_1 = \{i \mid 1 \leq i \leq l, i \in B \text{ and } i + l \in B\}, \quad B_1^* = \{i + l \mid i \in B_1\}, \\
B_2 = \{i \mid 1 \leq i \leq l, i \in B \text{ and } i + l \notin B\}, \\
B_3^* = \{i + l \mid 1 \leq i \leq l, i \notin B \text{ and } i + l \in B\}, \quad B_3 = \{i \mid i + l \in B_3^*\}.
\]

Then
\[
B = B_1 \cup B_1^* \cup B_2 \cup B_3^*.
\]

Thus
\[
s^T_B \bar{Q} B B^T s_B = s^T_B - s^T_B^* \begin{bmatrix} Q_{B_1B_1} & Q_{B_1B_2} & -Q_{B_1B_3} \\ Q_{B_2B_1} & Q_{B_2B_2} & -Q_{B_2B_3} \\ -Q_{B_3B_1} & -Q_{B_3B_2} & Q_{B_3B_3} \end{bmatrix} \begin{bmatrix} s_{B_1} - s_{B_1^*} \\ s_{B_2} \\ s_{B_3^*} \end{bmatrix}.
\]

Since
\[
\begin{bmatrix} Q_{B_1B_1} & Q_{B_1B_2} & -Q_{B_1B_3} \\ Q_{B_2B_1} & Q_{B_2B_2} & -Q_{B_2B_3} \\ -Q_{B_3B_1} & -Q_{B_3B_2} & Q_{B_3B_3} \end{bmatrix}
\]
has the same eigenvalues as
\[
\begin{bmatrix} Q_{B_1B_1} & Q_{B_1B_2} & Q_{B_1B_3} \\ Q_{B_2B_1} & Q_{B_2B_2} & Q_{B_2B_3} \\ Q_{B_3B_1} & Q_{B_3B_2} & Q_{B_3B_3} \end{bmatrix},
\]
which is
a square sub-matrix of $Q$, and $|B_1 \cup B_2 \cup B_3| \leq q$, we have
\[
-\frac{1}{2} s_B^T Q B B s_B 
\leq -\frac{\sigma}{2} \| \begin{bmatrix} s_{B_1} - s_{B_1}^* \\ s_{B_2} \\ s_{B_3}^* \end{bmatrix} \|^2 
\leq -\frac{\sigma}{2} \| \begin{bmatrix} s_{B_1}^* \\ s_{B_2}^* \\ s_{B_3} \end{bmatrix} \|^2 
= -\frac{\sigma}{2} \| s_B \|^2,
\]
where $\sigma = \min_I (\min(\text{eig}(Q_{II})))$, and $I$ is any subset of $\{1, \ldots, l\}$ with $|I| \leq q$. Note that $\|s_{B_1} - s_{B_1}^*\|^2 \geq \| \begin{bmatrix} s_{B_2} \\ s_{B_3}^* \end{bmatrix} \|^2$ is because of the following reasons: Since $\alpha_{i}^{k+1} (\alpha^*)_{i}^{k+1} = 0$, we consider two situations:

1. $\alpha_{i}^{k+1} = 0$: Then if $\alpha_{i}^{k} = 0$, $s_i = 0$ so $s_i s_{i+l} \leq 0$. On the other hand, if $\alpha_{i}^{k} > 0$, $(\alpha^*)_{i}^{k} = 0$. Hence $s_i \leq 0$ and $s_{i+l} \geq 0$ imply $s_i s_{i+l} \leq 0$.

2. $(\alpha^*)_{i}^{k+1} = 0$: Similarly, $s_i s_{i+l} \leq 0$.

Therefore, we have $-s_{B_1}^T s_{B_1}^* \geq 0$ so
\[
\|s_{B_1} - s_{B_1}^*\|^2 = \| \begin{bmatrix} s_{B_1}^* \\ s_{B_1} \end{bmatrix} \|^2 - 2 s_{B_1}^T s_{B_1}^* \geq \| \begin{bmatrix} s_{B_1} \\ s_{B_1}^* \end{bmatrix} \|^2.
\]

Recent implementation using Algorithms I.1 and I.2 with $q = 2$ for (40) are LIBSVM [5] and SVM Torch [6].

Note that it may be possible to extend convergence results in this section for algorithms used in [13], [15], [26] but here we will not get into details.

We then briefly discuss two other SVM formulations: one-class SVM and $\nu$-SVM. For one-class SVM [24], the formulation is already in the form of (37). For $\nu$-SVM [25], it has two linear constraints so is not covered by the algorithm and proof here.

The last formulation which we will consider is the support vector classification with quadratic penalty functions [7]. The required optimization problem is as follows:
\[
\min \frac{1}{2} \alpha^T (Q + I/C) \alpha - e^T \alpha \\
0 \leq \alpha_i, i = 1, \ldots, l, \\
y^T \alpha = 0,
\]
(43)
where \( I \) is the identity matrix.

As the upper bound of \( \alpha \) is \( \infty \) so the feasible region of (43) seems to be unbounded. This is a difficulty since we need the bounded property in order to have convergent subsequences.

For any feasible point \( \alpha \) of (43), since \( Q \) is positive semidefinite,

\[
\frac{1}{2C} \alpha^T \alpha - e^T \alpha \leq \frac{-1}{2} \alpha^T Q \alpha \leq 0.
\]

Therefore,

\[
\frac{1}{2C} \sum_{i=1}^{l} (\alpha_i - C)^2 \leq \frac{l}{2C} \cdot C^2 = \frac{lC}{2}
\]

implies that

\[
\alpha_i - C \leq \sqrt{lC}.
\]

Thus all feasible \( \alpha \) are in fact in a compact set. Then the convergence proof follows.

VI. Conclusions and Discussions

In this section we give some notes about the convergence proof. The property that (2) is exactly solved is used both in Lemmas IV.2 and IV.4. This confirms the conjectures in Section I where we think that an optimal solution of (2) makes a difference from the original Zoutendijk’s method.

It is unfortunate that we need Assumption IV.1 for the proof. We hope that this gap can be filled sometime in the future.

The convergence proof also suggests a possible way to improve the implementation. In final iterations, as the order of sorting \( y_i \nabla f(\alpha^*)_i, i = 1, \ldots, l \) is about fixed, it might be possible to consider fewer elements on the working set selection.

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