A Parallel Incremental Extreme SVM Classifier

Qing He

Institute of Computing Technology, Chinese Academy of Sciences University of Chinese Academy of Sciences

Outline

- SLFN and ELM
- Extreme SVM
- Parallel Extreme SVM
- Incremental Extreme SVM
- Conclusion

SLFN

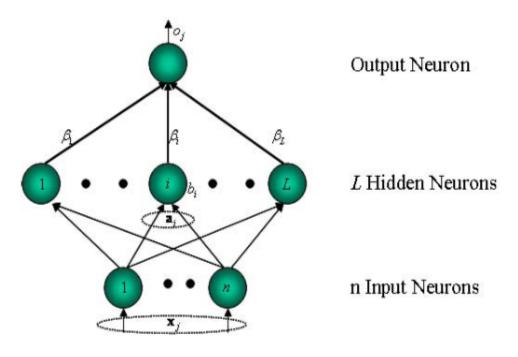


Figure 1: Feedforward Network Architecture: additive hidden nodes

Output of hidden nodes

$$G(\mathbf{a}_i, b_i, \mathbf{x}) = g(\mathbf{a}_i \cdot \mathbf{x} + b_i) \tag{1}$$

a_i: the weight vector connecting the *i*th hidden node and the input nodes.

b_i: the threshold of the *i*th hidden node.

Output of SLFNs

$$f_L(\mathbf{x}) = \sum_{i=1}^{L} \beta_i G(\mathbf{a}_i, b_i, \mathbf{x})$$
 (2)

 β_i : the weight vector connecting the *i*th hidden node and the output nodes.

SLFN

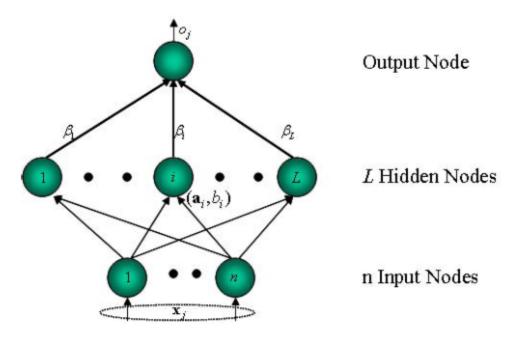


Figure 10: Feedforward Network Architecture

Mathematical Model

For N arbitrary distinct samples $(\mathbf{x}_i, \mathbf{t}_i) \in \mathbf{R}^n \times \mathbf{R}^m$, standard SLFNs with L hidden nodes and activation function g(x) are mathematically modeled as

$$\sum_{i=1}^{L} \beta_i G(\mathbf{a}_i, b_i, \mathbf{x}_j) = \mathbf{t}_j, \quad j = 1, \cdots, N$$
(7)

- a_i: the input weight vector connecting the ith hidden node and the input nodes or the center of the ith hidden node.
- β_i : the weight vector connecting the *i*th hidden node and the output node.
- b_i: the threshold or impact factor of the ith hidden node.

SLFN

Mathematical Model

 $\sum_{i=1}^{L} \beta_i G(\mathbf{a}_i, b_i, \mathbf{x}_i) = \mathbf{t}_i, j = 1, \cdots, N$ is equivalent to $\mathbf{H}\beta = \mathbf{T}$, where

$$H(a_{1}, \dots, a_{L}, b_{1}, \dots, b_{L}, x_{1}, \dots, x_{N})$$

$$= \begin{bmatrix} G(a_{1}, b_{1}, x_{1}) & \cdots & G(a_{L}, b_{L}, x_{1}) \\ \vdots & & \vdots \\ G(a_{1}, b_{1}, x_{N}) & \cdots & G(a_{L}, b_{L}, x_{N}) \end{bmatrix}_{N \times L}$$
(8)

$$\beta = \begin{bmatrix} \beta_1^T \\ \vdots \\ \beta_L^T \end{bmatrix}_{L \times m} \text{ and } \mathbf{T} = \begin{bmatrix} \mathbf{t}_1^T \\ \vdots \\ \mathbf{t}_N^T \end{bmatrix}_{N \times m}$$
 (9)

H is called the hidden layer output matrix of the neural network; the *i*th column of **H** is the output of the *i*th hidden node with respect to inputs x_1, x_2, \dots, x_N .

Extreme Learning Machine (ELM)

Three-Step Learning Model

Given a training set $\aleph = \{(\mathbf{x}_i, \mathbf{t}_i) | \mathbf{x}_i \in \mathbf{R}^n, \mathbf{t}_i \in \mathbf{R}^m, i = 1, \dots, N\}$, activation function g, and the number of hidden nodes L,

- Assign randomly input weight vectors or centers \mathbf{a}_i and hidden node bias or impact factor b_i , $i = 1, \dots, L$.
- Calculate the hidden layer output matrix H.
- **3** Calculate the output weight β : $\beta = \mathbf{H}^{\dagger}\mathbf{T}$.

where H[†] is the Moore-Penrose generalized inverse of hidden layer output matrix H.

Remarks on ELM

- It can be proved that $\hat{\beta} = \mathbf{H}^{\dagger}\mathbf{T}$ is the minimum norm least-squares solution of $\mathbf{H}\beta = \mathbf{T}$, which is unique.
- Minimum training error.

$$\|\mathbf{H}\hat{\beta} - \mathbf{T}\| = \|\mathbf{H}\mathbf{H}^{\dagger}\mathbf{T} - \mathbf{T}\| = \min_{\beta} \|\mathbf{H}\beta - \mathbf{T}\|.$$

This is because the solution is one of the leastsquares solutions

Smallest norm of weights. The special solution has the smallest norm among all the least-squares solutions.

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Linear ESVM

- Consider the 2-class classification problem of classifying m points in \mathbb{R}^n , represented by the $m \times n$ matrix A. A $m \times m$ diagonal matrix D with +1 or -1 along its diagonal specifies the membership of class A+ or class A- of each point A_i .
- For this problem, ESVM with a linear kernel, which has the same form as the linear PSVM, is given by the following quadratic program with parameter *v* > 0 and linear equality constraint (*y* is the slack variable):

$$\min_{(w,r,y)\in R^{n+1+m}} \frac{\nu}{2} ||y||^2 + \frac{1}{2} || \begin{bmatrix} w \\ r \end{bmatrix} ||^2$$
s.t.
$$D(Aw - er) + y = e$$

Nonlinear ESVM

- Suppose $\Phi(x)$ is the hidden layer output vector of an input vector \mathcal{X} in ELM algorithm. For the $m \times n$ matrix A, $\Phi(A)$ is defined as $\Phi(A) = [\Phi(A'_1), \dots, \Phi(A'_m)]'$
- Then, the nonlinear ESVM can be formulated to be the following quadratic program:

$$\min_{(w,r,y)\in R^{\tilde{n}+1+m}} \frac{\nu}{2} ||y||^2 + \frac{1}{2} || \begin{bmatrix} w \\ r \end{bmatrix} ||^2$$
s.t. $D(\Phi(A)w - er) + y = e$

where \widetilde{n} is the dimension of the feature space, i.e. the number of hidden nodes in ELM

Nonlinear ESVM

Substituting y by its explicit expression, we can get the following unconstrained minimization problem:

$$\min_{(w,r)\in R^{\widetilde{n}+1}} \frac{\nu}{2} \|D(\Phi(A)w - er) - e\|^2 + \frac{1}{2} \| \begin{bmatrix} w \\ r \end{bmatrix} \|^2$$

By setting the gradient with respect to w and r to zero we can obtain the solution for w and r in terms of problem data:

$$\begin{bmatrix} w \\ r \end{bmatrix} = (\frac{I}{\nu} + E_{\Phi}' E_{\Phi})^{-1} E_{\Phi}' De$$

where
$$E_{\Phi} = [\Phi(A) - e] \in \mathbb{R}^{m \times (\tilde{n}+1)}$$

Problems for ESVM

- For large-scale data, memory on single-processor limits the performance of extreme SVM.
 - Parallel Extreme SVM

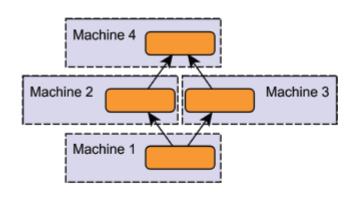
- For online settings, how can extreme SVM deal with new data?
 - Incremental Extreme SVM

Outline

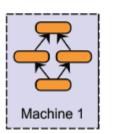
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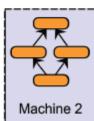
Parallel Strategy

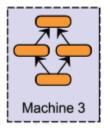
Model Parallelism

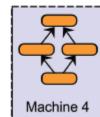


Data Parallelism

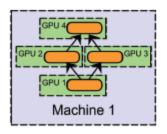


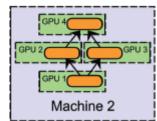


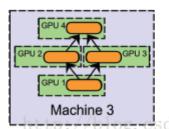


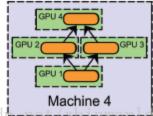


Model and Data Parallelism









Parallel ESVM

Our purpose is to parallel execute the computation for w and r

$$\begin{bmatrix} w \\ r \end{bmatrix} = (\frac{I}{\nu} + E_{\Phi}' E_{\Phi})^{-1} E_{\Phi}' De$$

- First, for a given $A \in R^{m \times n}$, to compute $\Phi(A) = [\Phi(A'_1), \dots, \Phi(A'_m)]' \in R^{m \times \tilde{n}}$, we can compute each $\Phi(A'_i)$, $i = 1, \dots, m$, in each parallel computing unit in principle
- To compute $E_{\Phi}'E_{\Phi}$, $E_{\Phi}'De$, consider the following decomposition:

$$E_{\Phi}'E_{\Phi} = [\alpha_1, \dots, \alpha_m][\alpha_1, \dots, \alpha_m]' = \sum_{i=1}^m \alpha_i \alpha_i'$$

$$E_{\Phi}'De = [\alpha_1, \dots, \alpha_m][d_1, \dots, d_m]' = \sum_{i=1}^m d_i \alpha_i$$

where α_i , $i=1,\ldots,m$, is an $\widetilde{n}+1$ dimensional column vector, and d_i , $i=1,\ldots,m$, is the class label of point A_i

Parallel ESVM

- In principle, we can compute each $\alpha_i \alpha_i'$, and $d_i \alpha_i$, i = 1, ..., m, in each parallel computing unit
- Since $\frac{I}{\nu} + E_{\Phi}' E_{\Phi}$ is a $(\widetilde{n} + 1) \times (\widetilde{n} + 1)$ matrix, The computation of $(\frac{I}{\nu} + E_{\Phi}' E_{\Phi})^{-1}$ is of order $(\widetilde{n} + 1)^3$; The computation of $(\frac{I}{\nu} + E_{\Phi}' E_{\Phi})^{-1} E_{\Phi}' De$ is of order $(\widetilde{n} + 1)^2$, where \widetilde{n} can be typically very small and is independent of the number of the training points m, so we serially execute them
- Algorithm 1 gives an explicit statement of our parallel ESVM algorithm. To make the algorithm clear, we assume that we have enough parallel computing units, and assign each input data point to a parallel computing unit

Parallel ESVM

Algorithm 1 Parallel ESVM classifier (PESVM)

Given an input dataset of m data points in R^n represented by the $m \times n$ matrix A and a diagonal matrix D of ± 1 labels denoting the class of each row of A, we generate the parallel ESVM classifier as follows:

- 1. Execute the following procedure for point A_i on the *i*-th parallel computing unit, where $i = 1, \ldots, m$:
 - (a) Compute $\Phi(A_i')$;
 - (b) Define $\alpha_i = \begin{bmatrix} \Phi(A_i') \\ -1 \end{bmatrix}$, and $d_i = D_{ii}$;
 - (c) Compute $\alpha_i \alpha_i'$ and $\bar{d}_i \alpha_i$;
- 2. Sum up $\alpha_i \alpha_i'$ and $d_i \alpha_i$ respectively, where i = 1, ..., m. And let $E_{\Phi}' E_{\Phi} = \sum_{i=1}^m \alpha_i \alpha_i', E_{\Phi}' De = \sum_{i=1}^m d_i \alpha_i$.
- 3. Compute the inversion $(\frac{I}{\nu} + E_{\Phi}' E_{\Phi})^{-1}$, for some positive value of ν . (Typically ν is chosen by means of a tuning set.)
- 4. Compute the solution $(\frac{I}{\nu} + E_{\Phi}' E_{\Phi})^{-1} E_{\Phi}' De$.

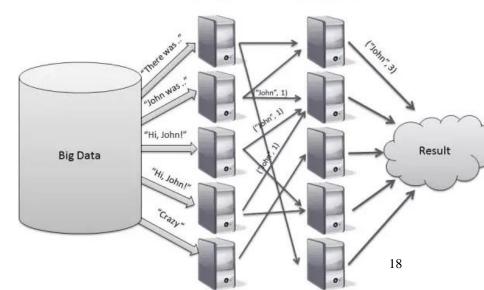
MapReduce





- Google
 - GFS, MapReduce, BigTable
 - Our abstraction is inspired by the map and reduce primitives present in Lisp and many other functional languages.
 - list = map(lambda x: x * x, [1,3,5,7,9])
 - sum = functools.reduce(lambda x,y: x+y, [1,3,5,7,9])
 MAP
 REDUCE
- Hadoop

Spark



- Parallel ESVM algorithm needs two Map/Reduce jobs. One for training (to be concrete, for getting $E_{\Phi}'E_{\Phi}$, $E_{\Phi}'De$), and the other for testing.
- In the training job, the map function first maps the input vector into a feature space by a random SLFN, then constructs α_i and conducts the vector products of $\alpha_i \alpha_i'$ and $\alpha_i d_i$. To decrease the cost of network communication, we define two global variables for each map task to accumulate each $\alpha_i \alpha_i'$ and $\alpha_i d_i$ within the same map task. We will collect these two global variables as the outputs in the close function of each map task (The close function executes when one map task is to be completed).

Algorithm 2 is the pseudocode of the map function of the training job

Algorithm 2 Train_map(key, value)

Input: Global variables localmatrix and localvector which are initialized with zeros, the offset key, the sample value

Output: Global variables local matrix and local vector which have accumulated $\alpha_i \alpha'_i$ and $\alpha_i d_i$ respectively

- 1. Construct the sample *instance* and the class label d_i from value;
- 2. Compute $\Phi(instance)$ where $\Phi(\cdot)$ is defined as (3);
- 3. Construct α_i as $\alpha_i = [\Phi(instance)' 1]'$;
- Compute α_iα'_i;
- 5. $localmatrix \leftarrow localmatrix + \alpha_i \alpha_i';$
- Compute α_id_i, where d_i is the class label of instance;
- 7. $localvector \leftarrow localvector + \alpha_i d_i$;

Algorithm 3. Train map close ().

Input: Global variables *localmatrix* and *localvector* which respectively represent the sum of local matrixes and local vectors within the same map task

Output: Global variables *localmatrix* and *localvector* which have accumulated $\alpha_i \alpha_i'$ and $\alpha_i d_i$ respectively < key', value' > pairs, where key' is a number, value' is a string comprises the values of a vector

- 1. for each p = 1, ..., the number of rows of *localmatrix* do
 - Take p as key';
- Construct value as a string, which comprises the values of localmatrix's pth row;
 - Emit < key', value' > pair;
 - 2. end for
 - 3. Take the number of rows of *localmatrix* + 1 as *key*;
- 4. Construct *value'* as a string, which comprises the values of *localvector*;
 - 5. Emit $\langle key \rangle$, value' > pair;

Finally the reduce function sums the outputs of each map task, which is illustrated in Algorithm 4:

Algorithm 4 Train_reduce(key, V)

Input: key is the row number of local matrix or the number of rows of local matrix + 1, V is the list of strings from different host, each of which comprises a local matrix's key-th row or a local vectorOutput: $\langle key', value' \rangle$ pairs, where key' is a row number, value' is a string which comprises the values of a vector, which represents the key'-th row of $E_{\Phi}'E_{\Phi}$, or $E_{\Phi}'De$

- Initialize one array temparray with zeros, to record the sum of the vectors in the list V;
- 2. while V.hasNext() do
 - Construct the vector vector from V.next();
 - temparray ← temparray + vector;
- 3. end while
- 4. Take key as key';
- Construct value' as a string, which comprises the values of temparray;
- 6. Emit $\langle key', value' \rangle$ pair;

 Algorithm 5 gives the pseudocode of the map function of the testing job. Reduce function is omitted.

Algorithm 5 Test_map(key, value)

Input: the offset key, the sample value, global variables w and r which is solution

Output: $\langle key', value' \rangle$ pairs, where key' is the flag of correct classification or wrong classification and value' is a string which comprises the sample information and the classification result

- Construct the sample instance and its class label classlabel from value;
- 2. Compute $\Phi(instance)$ where $\Phi(\cdot)$ is defined as (3);
- 3. Compute $\Phi(instance)'w r$;
- 4. if $\Phi(instance)'w r \ge 0$ then predictclass = A+; else

predictclass = A-;

5. if classlabel == predictclass then

```
key' = correct;
```

else

$$key' = wrong;$$

- Construct value' as a string, which comprises value and predictclass;
- 7. Emit $\langle key', value' \rangle$ pair;

Implementation based on Spark

Experimental settings:

- Spark version: Spark 2.2.0
- Java version: JDK1.8.0_112
- Servers: 3 CentOS servers, each with 48 cores and 128G memory

Dataset:

- Feature dimension: 64
- Size of training set: 79,999,211
- Size of testing set: 20,000,789

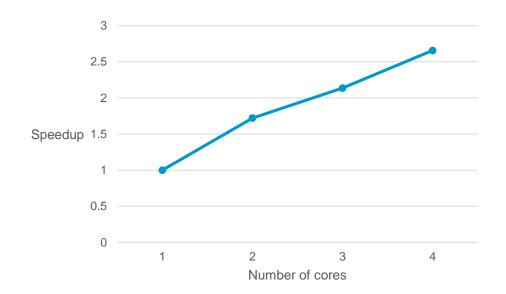
Classification Metric

Class	-1	1	
Precision	82.51%	87.65%	
Recall	88.56%	81.22%	
F1	85.43% 84.32%		
Accuracy	84.89%		

Speedup

Speedup =
$$\frac{\text{Execution time on 1 core}}{\text{Execution time on } m \text{ cores}}$$

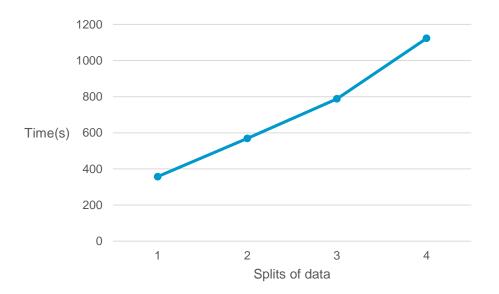
Number of cores	1	2	3	4
Execution Time(s)	1416	823	663	533
Speedup	1	1.72	2.13	2.65



Sizeup

$$Sizeup = \frac{Execution time for m splits of data}{Execution time for 1 split of data}$$

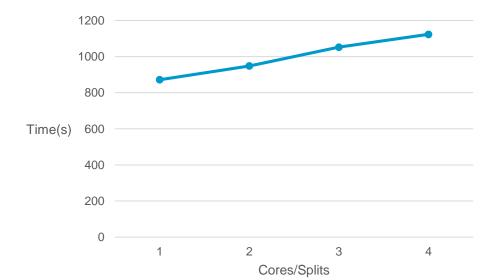
Splits of data (50M)	1	2	3	4
Execution Time(s)	357	569	788	1123
Sizeup	1	1.59	2.21	3.15



Scaleup

Scaleup = $\frac{\text{Execution time on 1 core for 1 split of data}}{\text{Execution time on } m \text{ cores for m splits of data}}$

Number of cores	1	2	3	4
Splits of data (50M)	1	2	3	4
Execution Time(s)	872	948	1052	1123



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Incremental ESVM

- Now we describe an incremental learning algorithm for ESVM that is capable of adding new data to generate an appropriately altered classifier.
- Assume that the current classifier is based on an input dataset, $A^1 \in \mathbb{R}^{m^1 \times n}$, and a corresponding $m^1 \times m^1$ diagonal matrix D^1 of ± 1 , and the classifier has the following form:

$$\begin{bmatrix} w \\ r \end{bmatrix} = (\frac{I}{\nu} + (E_{\Phi}^{1})' E_{\Phi}^{1})^{-1} (E_{\Phi}^{1})' D^{1} e$$

where
$$E_{\Phi}^1 = [\Phi(A^1) - e] \in \mathbb{R}^{m^1 \times (\tilde{n}+1)}$$

Suppose that a new set of training data points represented by the new matrix $A^2 \in \mathbb{R}^{m^2 \times n}$ needs to be added, for which we

Incremental ESVM

- correspondingly have $E_\Phi^2=[\Phi(A^2)-e]\in R^{m^2 imes(\widetilde n+1)}$, and an $m^2 imes m^2$ diagonal matrix D^2 of ± 1
- The classifier can be updated to reflect the addition of the new training dada as follows:

$$\begin{bmatrix} w \\ r \end{bmatrix} = \left(\frac{I}{\nu} + [(E_{\Phi}^{1})' (E_{\Phi}^{2})'] \begin{bmatrix} E_{\Phi}^{1} \\ E_{\Phi}^{2} \end{bmatrix} \right)^{-1} [(E_{\Phi}^{1})' (E_{\Phi}^{2})'] \begin{bmatrix} D^{1} O \\ O D^{2} \end{bmatrix} e$$

$$= \left(\frac{I}{\nu} + (E_{\Phi}^{1})' E_{\Phi}^{1} + (E_{\Phi}^{2})' E_{\Phi}^{2} \right)^{-1} ((E_{\Phi}^{1})' D^{1} e + (E_{\Phi}^{2})' D^{2} e)$$

We summarize our incremental ESVM algorithm as Algorithm 6

Incremental ESVM

Algorithm 6 Incremental ESVM classifier (IESVM)

Given a current classifier that is based on an input dataset of m^1 data points in R^n represented by the $m^1 \times n$ matrix A^1 and a diagonal matrix D of ± 1 labels denoting the class of each row of A^1 , we generate an incremental nonlinear classifier by adding new data represented by a new matrix $A^2 \in R^{m^2 \times n}$ and a corresponding diagonal matrix $D^2 \in R^{m^2 \times m^2}$ of ± 1 as follows:

- Compute Φ(A²) ∈ R^{m²×ñ} for the new data, which needs the random generated matrix W which
 is stored in the memory.
- 2. Define $E_{\Phi}^2 = [\Phi(A^2) e] \in \mathbb{R}^{m^2 \times (\widetilde{n}+1)}$.
- 3. Compute $(E_{\Phi}^2)'E_{\Phi}^2$ and $(E_{\Phi}^2)'D^2e$.
- Add the products obtained in step 3 to (E¹_Φ)'E¹_Φ and (E¹_Φ)'D¹e respectively, which are stored in the memory.
- 5. Compute the inversion of a $(\tilde{n}+1)\times(\tilde{n}+1)$ matrix and the product of this inversion and a $(\tilde{n}+1)$ dimensional vector.
- 6. The new solution is given by (11).

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Conclusion for ESVM

For large scale data, ESVM is easy to parallelize.

For online settings, ESVM could be learned incrementally.

Thank you!

 $Q. \mathcal{CIA}.$