Large-scale Linear Classification: Status and Challenges

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Talk at Criteo Machine Learning Workshop, November 8, 2017

Outline



- Optimization methods
- Multi-core linear classification
- Oistributed linear classification

5 Conclusions

Outline



- 2 Optimization methods
- 3 Multi-core linear classification
- Distributed linear classification

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Linear Classification

- Although many new and advanced techniques are available (e.g., deep learning), linear classifiers remain to be useful because of their simplicity
- We have fast training/prediction for large-scale data
- A large-scale optimization problem is solved
- The focus of this talk is on how to solve this optimization problem



The Software LIBLINEAR

• My talk will be very related to research done in developing the software LIBLINEAR for linear classification

www.csie.ntu.edu.tw/~cjlin/liblinear

 It is now one of the most used linear classification tools



Linear and Kernel Classification

Methods such as SVM and logistic regression are often used in two ways

• Kernel methods: data mapped to another space

$$\mathbf{x} \Rightarrow \phi(\mathbf{x})$$

 $\phi(\mathbf{x})^T \phi(\mathbf{y})$ easily calculated; no good control on $\phi(\cdot)$

 Feature engineering + linear classification: Directly use x without mapping. But x may have been carefully generated. Full control on x



Comparison Between Linear and Kernel

- For certain problems, accuracy by linear is as good as kernel
 - But training and testing are much faster
- Especially document classification
 Number of features (bag-of-words model) very large
 Large and sparse data
- Training millions of data in just a few seconds



Comparison Between Linear and Nonlinear (Training Time & Testing Accuracy)

	L	inear	RBF Kernel			
Data set	Time	Accuracy	Time	Accuracy		
MNIST38	0.1	96.82	38.1	99.70		
ijcnn1	1.6	91.81	26.8	98.69		
covtype_multiclass	1.4	76.37	46,695.8	96.11		
news20	1.1	96.95	383.2	96.90		
real-sim	0.3	97.44	938.3	97.82		
yahoo-japan	3.1	92.63	20,955.2	93.31		
webspam	25.7	93.35	15,681.8	99.26		
Size reasonably large: e.g., yahoo-japan: 140k instances and 830k features						

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and 830k features

Binary Linear Classification

- Training data $\{y_i, \boldsymbol{x}_i\}, \boldsymbol{x}_i \in R^n, i = 1, \dots, l, y_i = \pm 1$
- I: # of data, n: # of features

$$\min_{\boldsymbol{w}} \quad f(\boldsymbol{w}), \text{ where } f(\boldsymbol{w}) \equiv$$

$$C\sum_{i=1}^{l} \xi(\boldsymbol{w}; \boldsymbol{x}_{i}, y_{i}) + \begin{cases} \frac{1}{2} \boldsymbol{w}^{T} \boldsymbol{w} & \text{L2 regularization} \\ \|\boldsymbol{w}\|_{1} & \text{L1 regularization} \end{cases}$$

• $\xi(\boldsymbol{w}; \boldsymbol{x}, \boldsymbol{y})$: loss function: we hope $\boldsymbol{y}\boldsymbol{w}^T\boldsymbol{x} > 0$

• C: regularization parameter



Loss Functions

• Some commonly used loss functions.

$$\begin{aligned} \xi_{L1}(\boldsymbol{w};\boldsymbol{x},\boldsymbol{y}) &\equiv \max(0,1-\boldsymbol{y}\boldsymbol{w}^{T}\boldsymbol{x}), \quad (1) \\ \xi_{L2}(\boldsymbol{w};\boldsymbol{x},\boldsymbol{y}) &\equiv \max(0,1-\boldsymbol{y}\boldsymbol{w}^{T}\boldsymbol{x})^{2}, \quad (2) \\ \xi_{LR}(\boldsymbol{w};\boldsymbol{x},\boldsymbol{y}) &\equiv \log(1+e^{-\boldsymbol{y}\boldsymbol{w}^{T}\boldsymbol{x}}). \quad (3) \end{aligned}$$

- SVM (Boser et al., 1992; Cortes and Vapnik, 1995): (1)-(2)
- Logistic regression (LR): (3)

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 - 4 Distributed linear classification

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Optimization Methods

 A difference between linear and kernel is that for kernel, optimization must be over a variable α (usually through the dual problem) where

$$\mathbf{w} = \sum_{i=1}^{l} \alpha_i \phi(\mathbf{x}_i)$$

We cannot minimize over w, which may be infinite dimensional

• However, for linear, minimizing over w or α is ok



Optimization Methods (Cont'd)

Unconstrained optimization methods can be categorized to

- Low-order methods: quickly get a model, but slow final convergence
- High-order methods: more robust and useful for ill-conditioned situations

We will show both types of optimization methods are useful for linear classification

Further, to handle large problems, the algorithms must take problem structure into account

Let's discuss a low-order method (coordinate descent) in detail

Coordinate Descent

• We consider L1-loss and the dual SVM problem

$$\begin{array}{ll} \min_{\boldsymbol{\alpha}} & f(\boldsymbol{\alpha}) \\ \text{subject to} & 0 \leq \alpha_i \leq C, \forall i, \end{array}$$

$$f(\alpha) \equiv rac{1}{2} lpha^T Q lpha - oldsymbol{e}^T lpha$$

and

$$Q_{ij} = y_i y_j \boldsymbol{x}_i^T \boldsymbol{x}_j, \quad \boldsymbol{e} = [1, \dots, 1]^T$$

- \bullet We will apply coordinate descent (CD) methods
- The situation for L2 or LR loss is very similar

For current α, change α_i by fixing others
Let

$$\boldsymbol{e}_i = [0, \dots, 0, 1, 0, \dots, 0]^T$$

• The sub-problem is

$$\min_{d} f(\alpha + d\boldsymbol{e}_{i}) = \frac{1}{2}Q_{ii}d^{2} + \nabla_{i}f(\alpha)d + \text{constant}$$

subject to
$$0 \le \alpha_i + d \le C$$

• Without constraints

optimal
$$d=-rac{
abla_i f(oldsymbollpha)}{Q_{ii}}$$

• Now
$$0 \le \alpha_i + d \le C$$

$$\alpha_i \leftarrow \min\left(\max\left(\alpha_i - \frac{\nabla_i f(\boldsymbol{\alpha})}{Q_{ii}}, \mathbf{0}\right), C\right)$$

Note that

$$egin{aligned}
abla_i f(oldsymbollpha) &= (Qoldsymbollpha)_i - 1 = \sum_{j=1}^l oldsymbol Q_{ij} lpha_j - 1 \ &= \sum_{j=1}^l oldsymbol y_i oldsymbol y_j oldsymbol x_i^T oldsymbol x_j lpha_j - 1 \end{aligned}$$

• Expensive: O(In), I: # instances, n: features



• A trick in Hsieh et al. (2008) is to define and maintain

$$\mathbf{u}\equiv\sum_{j=1}^{l}y_{j}\alpha_{j}\mathbf{x}_{j},$$

• Easy gradient calculation: the cost is O(n)

$$\nabla_i f(\boldsymbol{\alpha}) = y_i \mathbf{u}^T \mathbf{x}_i - 1$$

 Note that this cannot be done for kernel as x_i is high dimensional



The procedure

• While α is not optimal (Outer iteration)

• For
$$i = 1, ..., l$$
 (Inner iteration)
(a) $\bar{\alpha}_i \leftarrow \alpha_i$
(b) $G = y_i \mathbf{u}^T \mathbf{x}_i - 1$
(c) $\alpha_i \leftarrow \min(\max(\alpha_i - G/Q_{ii}, 0), C)$
(d) If α_i needs to be changed
 $\mathbf{u} \leftarrow \mathbf{u} + (\alpha_i - \bar{\alpha}_i)y_i\mathbf{x}_i$

Maintaining \mathbf{u} also costs

O(n)

• Having

$$\mathbf{u} \equiv \sum_{j=1}^{l} y_j \alpha_j \mathbf{x}_j,$$
$$\nabla_i f(\boldsymbol{\alpha}) = y_i \mathbf{u}^T \mathbf{x}_i - 1$$

and

$$\bar{\alpha}_i$$
: old ; α_i : new
 $\mathbf{u} \leftarrow \mathbf{u} + (\alpha_i - \bar{\alpha}_i) y_i \mathbf{x}_i$.

is very essential

- This isn't the vanilla CD dated back to Hildreth (1957)
- We take the problem structure into account



Comparisons

L2-loss SVM is used

- DCDL2: Dual coordinate descent
- DCDL2-S: DCDL2 with shrinking
- PCD: Primal coordinate descent
- TRON: Trust region Newton method
- This result is from Hsieh et al. (2008) with C = 1



Optimization methods

Objective values (Time in Seconds)



Chih-Jen Lin (National Taiwan Univ.)

200

(IEI)

20

Low- versus High-order Methods

- We see low-order methods are efficient, but high-order methods are useful for difficult situations
- CD for dual
 - \$ time ./train -c 1 news20.scale
 - 2.528s
 - \$ time ./train -c 100 news20.scale
 28.589s
- Newton for primal
 - \$ time ./train -c 1 -s 2 news20.scale
 - 8.596s
 - \$ time ./train -c 100 -s 2 news20.scale
 11.088s



Training Median-sized Data: Status

- Basically a solved problem
- However, as data and memory continue to grow, new techniques are needed for large-scale sets.
- Two possible strategies are
 - Multi-core linear classification
 - Oistributed linear classification



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Multi-core Linear Classification

- Nowadays each CPU has several cores
- However, parallelizing algorithms to use multiple cores may not be that easy
- In fact, algorithms may need to be redesigned
- Since two years ago we have been working on multi-core LIBLINEAR



Multi-core Linear Classification (Cont'd)

- Three multi-core solvers have been released
 - Newton method for primal L2-regularized problem (Lee et al., 2015)
 - Coordinate descent method for dual L2-regularized problem (Chiang et al., 2016)
 - Coordinate descent method for primal L1-regularized problem (Zhuang et al., 2017)
- They are practically useful. For example, one user from USC thanked us because "a job (taking >30 hours using one core) now can finish within 5 hours"
- We will briefly discuss the 2nd and the 3rd



Multi-core CD for Dual

Recall the CD algorithm for dual is

• While α is not optimal (Outer iteration) For $i = 1, \ldots, l$ (Inner iteration) (a) $\bar{\alpha}_i \leftarrow \alpha_i$ (b) $G = y_i \mathbf{u}^T \mathbf{x}_i - 1$ (c) $\alpha_i \leftarrow \min(\max(\alpha_i - G/Q_{ii}, 0), C)$ (d) If α_i needs to be changed $\mathbf{u} \leftarrow \mathbf{u} + (\alpha_i - \bar{\alpha}_i) \mathbf{y}_i \mathbf{x}_i$



- The algorithm is inherently sequential
- Suppose

 $\alpha_{i'}$ is updated after α_i

Then $\alpha_{i'}$ must wait until the latest **u** is obtained

• The parallelization is difficult



- Asynchronous CD is possible (Hsieh et al., 2015), but may diverge
- We note that for a given set \bar{B}

$$abla_i f(\boldsymbol{w}) = \boldsymbol{w}^T \boldsymbol{x}_i, \forall i \in \bar{B}$$

can be calculated in parallel

• We then propose a framework



While α is not optimal
(a) Select a set B

(b) Calculate ∇_Bf(α) in parallel
(c) Select B ⊂ B with |B| ≪ |B|
(d) Sequentially update α_i, i ∈ B



• The selection of

$$B \subset \overline{B}$$
 with $|B| \ll |\overline{B}|$

is by $\nabla_{\bar{B}}f(w)$

• The idea is simple, but needs efforts to have a practical setting (details omitted)



Multi-core linear classification

Multi-core CD for Dual (Cont'd)



- Alg-4: the method in Chiang et al. (2016)
- Asynchronous CD (Hsieh et al., 2015)



Multi-core CD for L1 Regularization

- Currently, primal CD (Yuan et al., 2010) or its variants (Yuan et al., 2012) is the state-of-the-art for L1
- Each CD step involves one feature
- Some attempts of parallel CD for L1 include
 - Asynchronous CD (Bradley et al., 2011)
 - Block CD (Bian et al., 2013)
- These methods are not satisfactory for either
 - divergence issue, or
 - poor speedup



Multi-core CD for L1 Regularization (Cont'd)

- We struggled for years for find a solution
- Recently, in a work (Zhuang et al., 2017) we have an effective setting
- It's partially supported by Criteo Faculty Research Award
- Our idea is simple: direct parallelization of CD
- But wait.. This shouldn't work because each CD iteration is cheap



Direct Parallelization of CD

- Let's consider a simple setting to decide if one CD step should be parallelized or not
- if #non-zeros in an instance/feature \ge a threshold then

multi-core

else

single-core

Idea: a CD step is parallelized if there are enough operations



Direct Parallelization of CD (Cont'd)

• Speedup of CD for dual, L2 regularization

	#threads	_	Л	0
Data set			4	o
sparse	avazu-app	0.4	0.3	0.2
sets criteo	criteo	0.5	0.3	0.2
dense sets	epsilon_normalized	1.3	1.3	1.1
	splice_site.t.10%	1.8	2.8	4.1

- CD for dual: one instance at a time
- Threshold: 0 (sparse), 500 (dense)
- If 500 for sparse, no instance parallelized
- The speedup is poor



% of instances/features containing 50% and 80% # non-zeros

Data set	Instance		Feature		
avazu-app	50%	80%	0.2%	1%	
criteo	50%	80%	0.01%	0.2%	
kdd2010-a	40%	73%	0.03%	2%	
kdd2012	50%	80%	0.003%	0.5%	
rcv1_test	24%	54%	1%	5%	
splice_site.t.10%	50%	80%	9%	57%	
url_combined	44%	76%	0.002%	0.006%	
webspam	29%	55%	0.6%	2%	
yahoo-korea	20%	48%	0.07%	0.5%	

Features' non-zero distribution is extremely skewed Non-zeros are in few dense (and parallelizable) features



Speedup of CD for L1 Regularization

LR loss used	Naive			Block CD			Async. CD		
Data set	2	4	8	2	4	8	2	4	8
avazu-app	1.9	3.4	5.6	0.4	0.7	1.0	1.4	2.7	3.4
criteo	1.8	3.3	5.5	0.7	1.2	1.9	1.5	2.9	4.8
$epsilon_normalized$	2.0	4.0	7.9	x	х	х	1.3	2.1	х
HIGGS	2.0	3.9	7.5	0.7	0.8	0.9	1.0	1.3	х
kdd2010-a	1.7	2.4	3.1	0.8	1.4	2.4	1.5	2.7	4.8
kdd2012	1.9	2.8	3.9	0.2	0.4	0.6	2.1	4.7	7.0
rcv1_test	1.9	3.4	5.9	x	х	х	1.3	2.5	4.5
splice_site.t.10%	1.9	3.6	6.2	x	х	х	1.6	2.7	4.3
url_combined	2.0	3.5	6.2	0.5	0.9	1.3	1.0	1.7	1.7
webspam	1.8	3.2	4.8	0.1	0.3	0.5	1.4	2.5	4
vahoo-korea	1.9	3.5	5.9	0.2	0.3	0.5	1.3	2.4	4.4

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Distributed Linear Classification

- It's even more complicated than multi-core
- I don't have time to discuss this topic in detail, but let me share some lessons
- A big mistake was that we worked on distributed before multi-core



Distributed Linear Classification (Cont'd)

- A few years ago, big data was hot. So we extended a Newton solver in LIBLINEAR to MPI (Zhuang et al., 2015) and Spark (Lin et al., 2014)
- We were a bit ahead of time; Spark MLlib wasn't even available then
- Unfortunately, very few people use our code, especially the Spark one
- We moved to multi-core. Immediately, multi-core LIBLINEAR has many users



Distributed Linear Classification (Cont'd)

- Why we failed? Several possible reasons
- Not many people have big data??
- System issues are more important than we thought. At that time Spark wasn't easy to use and was being actively changed
- System configuration and application scenarios may significantly vary

An algorithm useful for systems with fast network speed may be useless for systems with slow communication



Distributed Linear Classification (Cont'd)

• Application dependency is stronger.

L2 and L1 regularization often give similar accuracy. On a single machine, we may not want to use L1 because training is more difficult and the smaller model size isn't that important

However, for distributed applications many have told me that they need $\ensuremath{\mathsf{L1}}$

• A lesson is that for people from academia, it's better to collaborate with industry for research on distributed machine learning



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Conclusions

- Linear classification is an old topic, but it remains to be useful for many applications
- Efficient training relies on designing optimization algorithms by incorporating the problem structure
- Many issues about multi-core and distributed linear classification still need to be studied

