

Supplementary Materials of “Limited-memory Common-directions Method for Distributed L1-regularized Linear Classification”

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I Introduction

In this document, we present additional details and more experimental results.

II Derivation of the Direction Used in LBFGS

By using only the information from the last m iterations, the definition of B_k in BFGS becomes the following in LBFGS.

$$(II.1) \quad B_k = V_{k-1}^T \cdots V_{k-m}^T B_0^k V_{k-m} \cdots V_{k-1} + \rho_{k-m} V_{k-1}^T \cdots V_{k-m+1}^T \mathbf{s}_{k-m} \mathbf{s}_{k-m}^T V_{k-m+1} \cdots V_{k-1} + \cdots + \rho_{k-1} \mathbf{s}_{k-1} \mathbf{s}_{k-1}^T.$$

Note that in BFGS, B_0^k is a fixed matrix, but in LBFGS, B_0^k can change with k , provided its eigenvalues are bounded in a positive interval over k . A common choice is

$$(II.2) \quad B_0^k = \frac{\mathbf{s}_{k-1}^T \mathbf{u}_{k-1}}{\mathbf{u}_{k-1}^T \mathbf{u}_{k-1}} I.$$

By expanding (II.1), \mathbf{d}_k can be efficiently obtained by $\mathcal{O}(m)$ vector operations as shown in Algorithm II. The overall procedure of LBFGS is summarized in Algorithm I.

III More Details of Limited-memory Common-directions Method

A sketch of the procedure for L1-regularized problems is in Algorithm III.

IV Line Search in Algorithms for L2-regularized Problems

Here we present the trick mentioned in Section 2.3 in the paper. At each line search iteration, we obtain $\mathbf{w}^T \mathbf{x}_i, \mathbf{d}^T \mathbf{x}_i, \forall i$ first, and then use $\mathcal{O}(l)$ cost to calculate

$$(\mathbf{w} + \alpha \mathbf{d})^T \mathbf{x}_i = \mathbf{w}^T \mathbf{x}_i + \alpha \mathbf{d}^T \mathbf{x}_i, \forall i.$$

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Algorithm I LBFGS.

- 1: Given \mathbf{w}_0 , integer $m > 0$, and $\beta, \gamma \in (0, 1)$
- 2: $\mathbf{w} \leftarrow \mathbf{w}_0$
- 3: **for** $k = 0, 1, 2, \dots$ **do**
- 4: Calculate $\nabla f(\mathbf{w})$
- 5: Calculate the new direction

$$\mathbf{d} \equiv -B \nabla f(\mathbf{w}) \approx -\nabla^2 f(\mathbf{w})^{-1} \nabla f(\mathbf{w})$$

by using information of the previous m iterations (Algorithm II)

- 6: Calculate $\nabla f(\mathbf{w})^T \mathbf{d}$
 - 7: $\alpha \leftarrow 1, \mathbf{w}^{\text{old}} \leftarrow \mathbf{w}$
 - 8: **while** true **do**
 - 9: $\mathbf{w} \leftarrow \mathbf{w}^{\text{old}} + \alpha \mathbf{d}$
 - 10: Calculate the objective value $f(\mathbf{w})$ in (2.2)
 - 11: **if** $f(\mathbf{w}) - f(\mathbf{w}^{\text{old}}) \leq \gamma \nabla f(\mathbf{w}^{\text{old}})^T (\mathbf{w} - \mathbf{w}^{\text{old}})$
 - 12: **break**
 - 13: $\alpha \leftarrow \alpha \beta$
 - 14: Update P with $\mathbf{w} - \mathbf{w}^{\text{old}}$ and $\nabla f(\mathbf{w}) - \nabla f(\mathbf{w}^{\text{old}})$
-

Because $\mathbf{w}^T \mathbf{x}_i$ can be obtained from the previous iteration, $\mathbf{d}^T \mathbf{x}_i$ is the only $\mathcal{O}(\#\text{nnz})$ operation needed. The line search cost is reduced from

$$\#\text{line-search steps} \times \mathcal{O}(\#\text{nnz})$$

to

$$1 \times \mathcal{O}(\#\text{nnz}) + \#\text{line-search steps} \times \mathcal{O}(l).$$

This trick is not applicable to L1-regularized problems because the new point is no longer $\mathbf{w} + \alpha \mathbf{d}$.

V More on the Distributed Implementation

V.1 Complexity. The distributed implementation as mentioned in Section 5 is shown in Algorithm VI. Then we discuss the complexity below.

$$\frac{(2 + \#\text{line-search steps}) \times \mathcal{O}(\#\text{nnz}) + \mathcal{O}(lm^2) + \mathcal{O}(mn)}{K} + \mathcal{O}(m^3)$$

VI More Experiments

The data sets used in this section are shown in Table (I).

Algorithm II LBFSG Two-loop recursion.

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1:  $\mathbf{q} \leftarrow -\nabla f(\mathbf{w})$ 
2: for  $i = k - 1, k - 2, \dots, k - m$  do
3:    $\alpha_i \leftarrow \mathbf{s}_i^T \mathbf{q} / \mathbf{s}_i^T \mathbf{u}_i$ 
4:    $\mathbf{q} \leftarrow \mathbf{q} - \alpha_i \mathbf{u}_i$ 
5:  $\mathbf{r} \leftarrow (\mathbf{s}_{k-1}^T \mathbf{u}_{k-1} / \mathbf{u}_{k-1}^T \mathbf{u}_{k-1}) \mathbf{q}$ 
6: for  $i = k - m, k - m + 1, \dots, k - 1$  do
7:    $\beta_i \leftarrow \mathbf{u}_i^T \mathbf{r} / \mathbf{s}_i^T \mathbf{u}_i$ 
8:    $\mathbf{r} \leftarrow \mathbf{r} + (\alpha_i - \beta_i) \mathbf{s}_i$ 
9:  $\mathbf{d} \leftarrow \mathbf{r}$ 

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Algorithm III Limited-memory common-directions method for L1-regularized problems.

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1: while true do
2:   Compute  $\nabla^P f(\mathbf{w})$  by (2.6)
3:   Solve the sub-problem (3.21)
4:   Let the direction be  $\mathbf{d} = P\mathbf{t}$ 
5:   for  $j = 1, \dots, n$  do
6:     Align  $d_j$  with  $-\nabla_j^P f(\mathbf{w})$  by (2.11)
7:    $\alpha \leftarrow 1, \mathbf{w}^{\text{old}} \leftarrow \mathbf{w}$ 
8:   while true do
9:     Calculate  $\mathbf{w}$  from  $\mathbf{w}^{\text{old}} + \alpha \mathbf{d}$  by (2.12)
10:    if  $f(\mathbf{w}) - f(\mathbf{w}^{\text{old}}) \leq \gamma \nabla^P f(\mathbf{w}^{\text{old}})^T (\mathbf{w} - \mathbf{w}^{\text{old}})$ 
11:      break
12:     $\alpha \leftarrow \alpha \beta$ 
13:   Update  $P$  and  $XP$ 

```

VI.1 More Results by Using Different C Values. In Figure (I), we present more results with

$$C = \{0.1C_{\text{Best}}, C_{\text{Best}}, 10C_{\text{Best}}\},$$

where C_{Best} is the value to achieve the highest cross validation accuracy. The results are similar to C_{Best} presented in Section 6, but we observe that NEWTON converges slowly in larger C cases.

VI.2 More Results on Distributed Experiments. In Figure (II), we present more data sets for the distributed experiments. All settings are the same as in Section 6.

Table (I): Data statistics.

Data set	#instances	#features	#nonzeros	C_{Best}
real-sim	72,309	20,958	3,709,083	16
rcv1_test	677,399	47,226	49,556,258	4
news20	19,996	1,355,191	9,097,916	1024
yahoojp	176,203	832,026	23,506,415	2
url	2,396,130	3,231,961	277,058,644	8
yahookr	460,554	3,052,939	156,436,656	4
epsilon	400,000	2,000	800,000,000	0.5
webspam	350,000	16,609,143	1,304,697,446	64
KDD2010-b	19,264,097	29,890,096	566,345,888	0.5
criteo	45,840,617	1,000,000	1,787,773,969	0.5
avazu-site	25,832,830	999,962	387,492,144	1
kdd2012	149,639,105	54,686,452	1,646,030,155	2

Algorithm IV A distributed implementation of OWLQN.

- 1: **for** $k = 0, 1, 2, \dots$ **do**
- 2: Compute $\nabla^P f(\mathbf{w})$ by (2.6) and

$$\nabla L(\mathbf{w}) = C \bigoplus_{r=1}^K (X_{J_r, :})^T \begin{bmatrix} \vdots \\ \xi^l(y_i \mathbf{w}^T \mathbf{x}_i) \\ \vdots \end{bmatrix}_{i \in J_r}.$$

- $\triangleright \mathcal{O}(\#\text{nnz}/K)$; $\mathcal{O}(n)$ comm.
- 3: Compute the search direction $\mathbf{d}_{\bar{J}_r}, r = 1, \dots, K$
by Algorithm V $\triangleright \mathcal{O}(nm/K)$; $\mathcal{O}(m)$ comm.
- 4: An *allgather* operation to let each node has

$$\mathbf{d} = \begin{bmatrix} \mathbf{d}_{\bar{J}_1} \\ \vdots \\ \mathbf{d}_{\bar{J}_K} \end{bmatrix}$$

$\triangleright \mathcal{O}(n/K)$ comm.

- 5: **for** $j = 1, \dots, n$ **do**
- 6: Align d_j with $-\nabla_{\bar{J}_j}^P f(\mathbf{w})$ by (2.11)
- 7: $\alpha \leftarrow 1, \mathbf{w}^{\text{old}} \leftarrow \mathbf{w}$
- 8: **while** true **do**
- 9: Calculate \mathbf{w} from $\mathbf{w}^{\text{old}} + \alpha \mathbf{d}$ by (2.12) and

$$f(\mathbf{w}) = \|\mathbf{w}\|_1 + C \bigoplus_{r=1}^K \sum_{i \in J_r} \xi(y_i \mathbf{w}^T \mathbf{x}_i)$$

$\triangleright \mathcal{O}(\#\text{nnz}/K)$; $\mathcal{O}(1)$ comm.

- 10: **if** $f(\mathbf{w}) - f(\mathbf{w}^{\text{old}}) \leq \gamma \nabla^P f(\mathbf{w}^{\text{old}})^T (\mathbf{w} - \mathbf{w}^{\text{old}})$
- 11: **break**
- 12: $\alpha \leftarrow \alpha \beta$
- 13: $\mathbf{s}_k \leftarrow \mathbf{w} - \mathbf{w}^{\text{old}}, \mathbf{u}_k \leftarrow \nabla f(\mathbf{w}) - \nabla f(\mathbf{w}^{\text{old}})$
- 14: Remove 1st column of S and U if needed and

$$S \leftarrow [S \quad \mathbf{s}_k], \quad U \leftarrow [U \quad \mathbf{u}_k]$$

- 15: $\rho_k \leftarrow \bigoplus_{r=1}^K (\mathbf{u}_k)_{\bar{J}_r}^T (\mathbf{s}_k)_{\bar{J}_r} \quad \triangleright \mathcal{O}(n/K)$; $\mathcal{O}(1)$ comm.
-

Algorithm V Distributed OWLQN Two-loop recursion

- 1: **if** $k = 0$ **return** $\mathbf{d}_{\bar{J}_r} \leftarrow -\nabla_{\bar{J}_r}^P f(\mathbf{w})$
- 2: **for** $r = 1, \dots, K$ **do in parallel**
- 3: $\mathbf{q}_{\bar{J}_r} \leftarrow -\nabla_{\bar{J}_r}^P f(\mathbf{w}) \quad \triangleright \mathcal{O}(n/K)$
- 4: **for** $i = k-1, k-2, \dots, k-m$ **do**
- 5: Calculate α_i by

$$\alpha_i \leftarrow \frac{\bigoplus_{r=1}^K (\mathbf{s}_i)_{\bar{J}_r}^T \mathbf{q}_{\bar{J}_r}}{\rho_i}$$

$\triangleright \mathcal{O}(n/K)$; $\mathcal{O}(1)$ comm.

- 6: **for** $r = 1, \dots, K$ **do in parallel**
- 7: $\mathbf{q}_{\bar{J}_r} \leftarrow \mathbf{q}_{\bar{J}_r} - \alpha_i (\mathbf{u}_i)_{\bar{J}_r} \quad \triangleright \mathcal{O}(n/K)$
- 8: Calculate

$$\mathbf{u}_{k-1}^T \mathbf{u}_{k-1} \leftarrow \bigoplus_{r=1}^K (\mathbf{u}_{k-1})_{\bar{J}_r}^T (\mathbf{u}_{k-1})_{\bar{J}_r}$$

$\triangleright \mathcal{O}(n/K)$; $\mathcal{O}(1)$ comm.

- 9: **for** $r = 1, \dots, K$ **do in parallel**
- 10: $\mathbf{r}_{\bar{J}_r} \leftarrow \frac{\rho_{k-1}}{\mathbf{u}_{k-1}^T \mathbf{u}_{k-1}} \mathbf{r}_{\bar{J}_r} \quad \triangleright \mathcal{O}(n/K)$
- 11: **for** $i = k-m, k-m+1, \dots, k-1$ **do**
- 12: Calculate β_i by

$$\beta_i \leftarrow \frac{\bigoplus_{r=1}^K (\mathbf{u}_i)_{\bar{J}_r}^T \mathbf{r}_{\bar{J}_r}}{\rho_i}$$

$\triangleright \mathcal{O}(n/K)$; $\mathcal{O}(1)$ comm.

- 13: **for** $r = 1, \dots, K$ **do in parallel**
 - 14: $\mathbf{r}_{\bar{J}_r} \leftarrow \mathbf{r}_{\bar{J}_r} + (\alpha_i - \beta_i) (\mathbf{s}_i)_{\bar{J}_r} \quad \triangleright \mathcal{O}(n/K)$
 - return** $\mathbf{d}_{\bar{J}_r} \leftarrow \mathbf{r}_{\bar{J}_r}, r = 1, \dots, K$
-

Algorithm VI Distributed limited-memory common-directions method.

- 1: **while** true **do**
- 2: Compute $\nabla^P f(\mathbf{w})$ by (2.6) and

$$\nabla L(\mathbf{w}) = C \bigoplus_{r=1}^K (X_{J_r,:})^T \begin{bmatrix} \vdots \\ \xi'(y_i; \mathbf{w}^T \mathbf{x}_i) \\ \vdots \end{bmatrix}_{i \in J_r}.$$

▷ $\mathcal{O}(\#\text{nnz}/K)$; $\mathcal{O}(n)$ comm.

- 3: Calculate

$$X_{J_r,:} \nabla^P f(\mathbf{w})$$

▷ $\mathcal{O}(\#\text{nnz}/K)$

- 4: Remove 1st column of P and U if needed and

$$\begin{aligned} P_{J_r,:} &\leftarrow [P_{J_r,:} \quad \nabla_{J_r}^P f(\mathbf{w})] \\ U_{J_r,:} &\leftarrow [U_{J_r,:} \quad X_{J_r,:} \nabla^P f(\mathbf{w})] \end{aligned}$$

- 5: Calculate

$$\begin{aligned} (XP)^T D_{\mathbf{w}}(XP) &= \bigoplus_{r=1}^K (U_{J_r,:})^T (D_{\mathbf{w}})_{J_r, J_r} U_{J_r,:} \\ &\quad \triangleright \mathcal{O}(lm^2/K), \mathcal{O}(m^2) \text{ comm.} \end{aligned}$$

$$\begin{aligned} -P^T \nabla^P f(\mathbf{w}) &= -\bigoplus_{r=1}^K (P_{\bar{J}_r,:})^T \nabla_{\bar{J}_r}^P f(\mathbf{w}) \\ &\quad \triangleright \mathcal{O}(mn/K); \mathcal{O}(m) \text{ comm.} \end{aligned}$$

- 6: Solve

$$\begin{aligned} ((XP)^T D_{\mathbf{w}}(XP)) \mathbf{t} &= -P^T \nabla^P f(\mathbf{w}) \\ &\quad \triangleright \mathcal{O}(m^3) \end{aligned}$$

- 7: Let the direction be

$$\begin{aligned} \mathbf{d} = P\mathbf{t} &= [P_{\bar{J}_1,:} \mathbf{t}, \dots, P_{\bar{J}_K,:} \mathbf{t}]^T \\ &\quad \triangleright \mathcal{O}(mn/K); \mathcal{O}(n/K) \text{ comm.} \end{aligned}$$

- 8: **for** $j = 1, \dots, n$ **do**

- 9: Align d_j with $-\nabla_j^P f(\mathbf{w})$ by (2.11)

- 10: $\alpha \leftarrow 1$, $\mathbf{w}^{\text{old}} \leftarrow \mathbf{w}$

- 11: **while** true **do**

- 12: Calculate \mathbf{w} from $\mathbf{w}^{\text{old}} + \alpha \mathbf{d}$ by (2.12) and

$$\begin{aligned} f(\mathbf{w}) &= \|\mathbf{w}\|_1 + C \bigoplus_{r=1}^K \sum_{i \in J_r} \xi(y_i; \mathbf{w}^T \mathbf{x}_i) \\ &\quad \triangleright \mathcal{O}(\#\text{nnz}/K); \mathcal{O}(1) \text{ comm.} \end{aligned}$$

- 13: **if** $f(\mathbf{w}) - f(\mathbf{w}^{\text{old}}) \leq \gamma \nabla^P f(\mathbf{w}^{\text{old}})^T (\mathbf{w} - \mathbf{w}^{\text{old}})$

- 14: **break**

- 15: $\alpha \leftarrow \alpha \beta$

- 16: Remove 1st column of P and U if needed and

$$\begin{aligned} P &\leftarrow [P \quad \mathbf{w} - \mathbf{w}^{\text{old}}] \\ U_{J_r,:} &\leftarrow [U_{J_r,:} \quad X_{J_r,:} (\mathbf{w} - \mathbf{w}^{\text{old}})] \end{aligned}$$

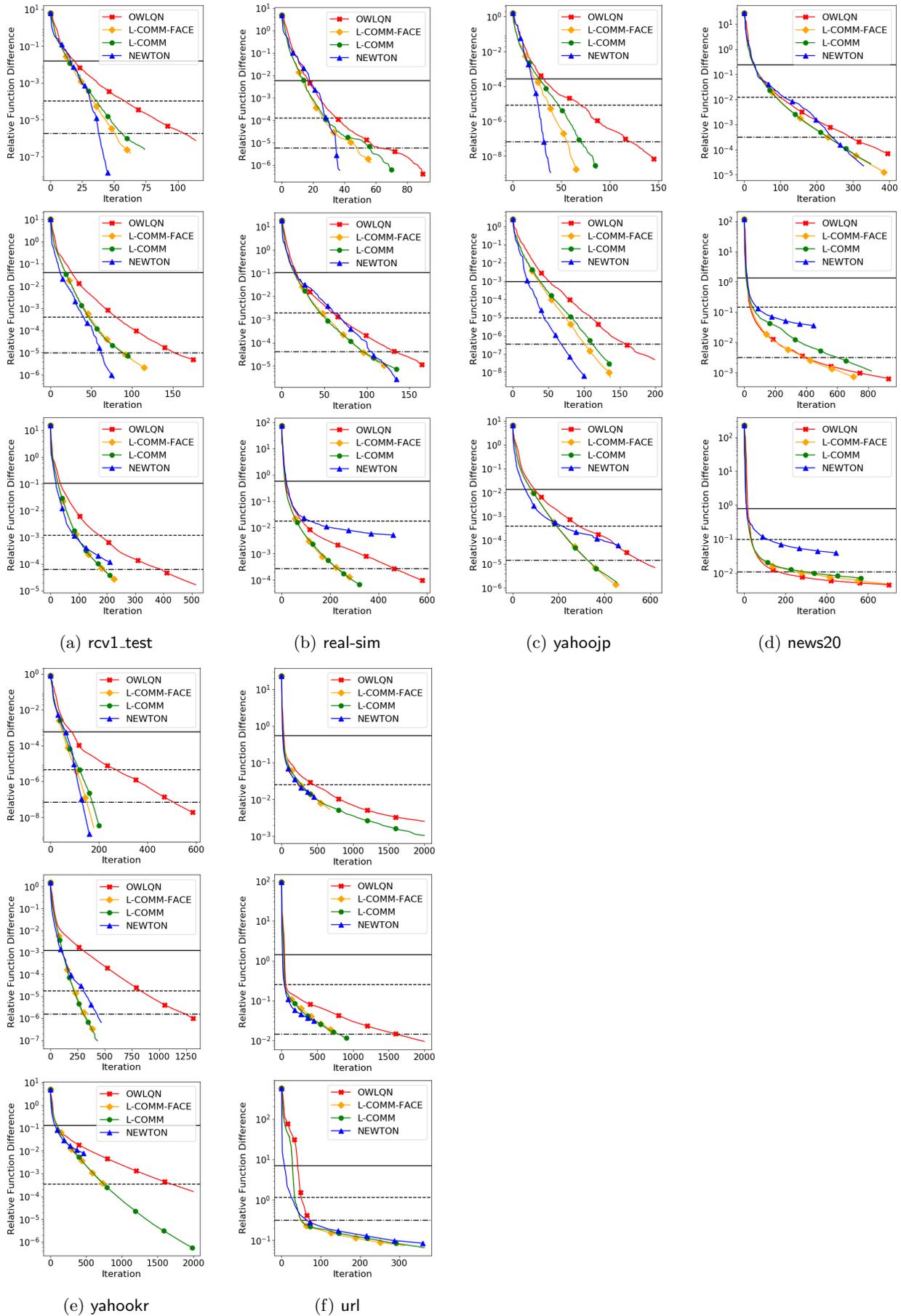


Figure (I): Comparison of different algorithms with $0.1C_{\text{Best}}$, C_{Best} , $10C_{\text{Best}}$, respectively from top to below for each data set. We show iteration versus the relative difference to the optimal value. Other settings are the same as in Figure 3

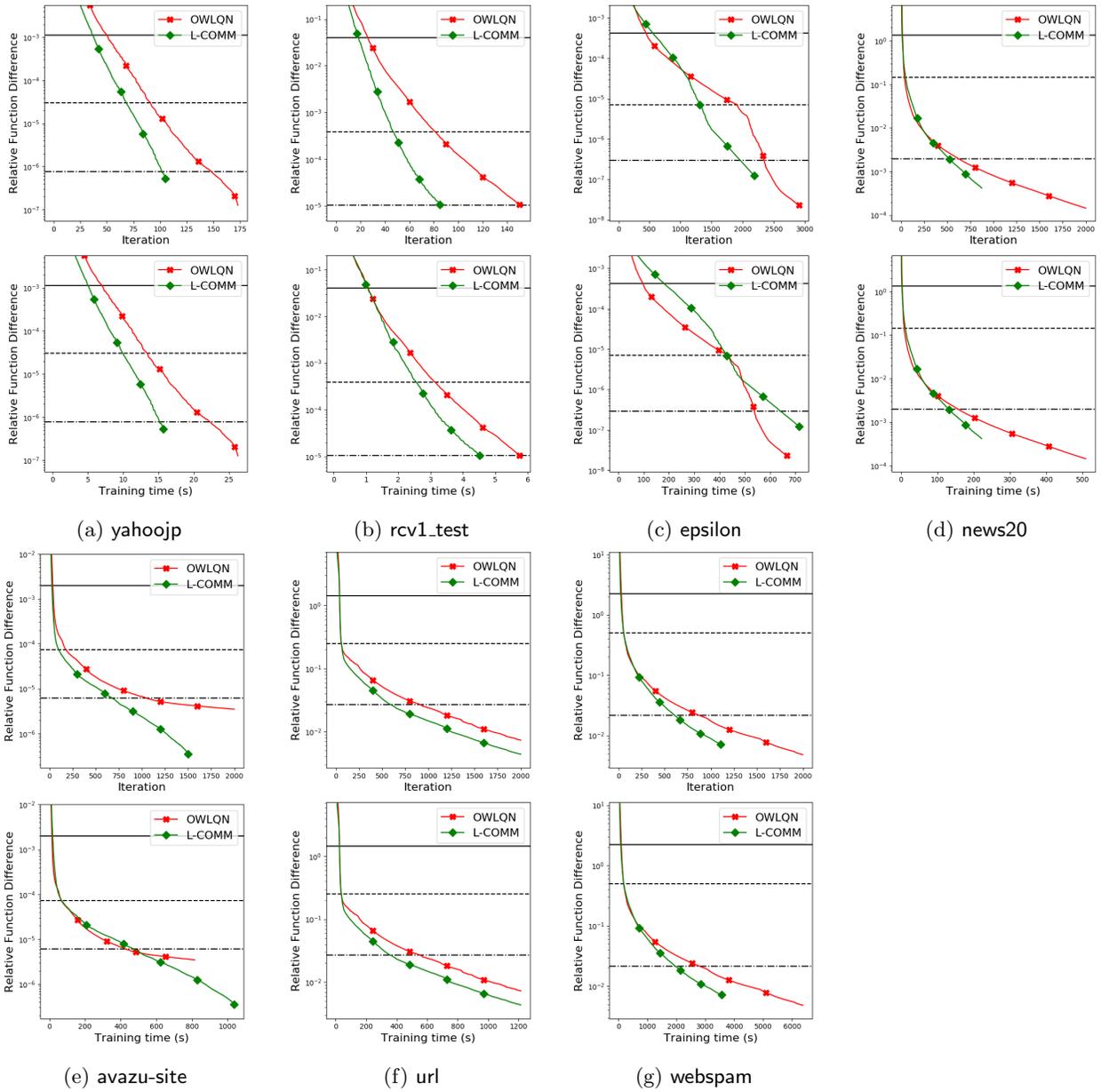


Figure (II): Comparison of different algorithms by using 32 nodes. Upper: iterations. Lower: running time in seconds. Other settings are the same as Figure 4.