

Weakly Supervised Open-set Domain Adaptation by Dual-domain Collaboration Supplementary Material

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1. Optimization Method

We define the total loss function as

$$f = Dist_C + \lambda_M Dist_M + \lambda_G G + \lambda_U U. \quad (1)$$

Then, our objective function can be concluded as follows:

$$(W_A^*, W_B^*) = \arg \min_{W_A, W_B} f(\mathcal{D}_A, \mathcal{D}_B, W_A, W_B). \quad (2)$$

The loss function f is not convex in the elements of W_A and W_B . Therefore we use the non-linear methods in [23] to optimize the transform matrices.

We jointly optimize W_A and W_B by concatenating them as a whole matrix $W \in \mathbb{R}^{d \times 2d}$. By optimizing W as a whole, we can ensure that both matrices are simultaneously updated. As for gradient $\partial f / \partial W$, we similarly concatenate $\partial f / \partial W_A$ and $\partial f / \partial W_B$ to form a single gradient value. After the optimization of W_A and W_B , we can easily obtain the projected features of \mathcal{D}_A and \mathcal{D}_B from $\mathcal{D}_A W_A$ and $\mathcal{D}_B W_B$.

To save space, we only give the gradient of each component in f for W_A . The gradient of W_B can be similarly formed. From Equation

$$f = Dist_C + \lambda_M Dist_M + \lambda_G G + \lambda_U U \quad (3)$$

we have

$$\frac{\partial f}{\partial W_A} = \frac{\partial Dist_C}{\partial W_A} + \lambda_M \frac{\partial Dist_M}{\partial W_A} + \lambda_G \frac{\partial G}{\partial W_A} + \lambda_U \frac{\partial U}{\partial W_A} \quad (4)$$

For the first three components, the gradient can be simply formulated as

$$\frac{\partial Dist_C}{\partial W_A} = \sum_{c \in \mathcal{C}_K} (P_A^c)^T (P_A^c W_A - P_B^c W_B), \quad (5)$$

$$\frac{\partial Dist_M}{\partial W_A} = (P_A)^T (P_A W_A - P_B W_B), \quad (6)$$

$$\frac{\partial G}{\partial W_A} = \sum_{c \in \mathcal{C}_L} \frac{1}{n_A^c} \sum_{x^i \in \mathbf{X}_A^c} (x^i - \bar{x}_A^c)^T (x^i - \bar{x}_A^c) W_A. \quad (7)$$

As for U , the gradient is composed of two parts

$$\frac{\partial U}{\partial W_A} = \frac{\partial U_A}{\partial W_A} + \frac{\partial U_B}{\partial W_A} \quad (8)$$

We first find indexes $(c, i) \in \mathcal{N}$ that triggers U_A , then,

$$\frac{\partial U_A}{\partial W_A} = \sum_{(c, i) \in \mathcal{N}} (x^i)^T (x_u W_{D'} - \bar{x}_A^c W_A), \quad (9)$$

where D' denotes the host domain of x_u . Then, we find the indexes $(c', i') \in \mathcal{N}$ that triggers U_B , where the nearest unknown neighbor sample x_u is from \mathcal{D}_A . Therefore, the gradient for U_B can be defined as:

$$\frac{\partial U_B}{\partial W_A} = \sum_{(c', i') \in \mathcal{N}} (x_u)^T (x_u W_A - x^{i'} W_B). \quad (10)$$

2. Time Complexity

We empirically compared the complexity of our method with top competitors on Intel E5-2650 CPU, with Nvidia GTX TITAN X GPU for the deep method. The time cost of each methods on different tasks are shown in Table 1. As shown, time complexity of CDA is comparable to other domain adaptation methods, while its accuracy outperformed all the compared methods.

3. Effect of Pseudo Labelling

We have compared this process with OSVM [9] and 1-vs-Set SVM [20]. We show the results for the first iteration of $A \leftrightarrow W$ in Table 3, including accuracies of different

Methods	A↔D	A↔W	W↔D	5↔6	7↔8
MMDT [8]	10.23	9.56	9.00	368.80	414.72
TCA [18]	2.04	1.92	0.59	25.41	39.04
OpenBP [19]	16109.38	11120.06	6061.15	6445.76	12559.27
CDA	7.28	8.03	4.72	112.05	169.10

Table 1. Time complexity of different methods (second).

kinds of samples, rates of unknown-class samples *mis-labeled* as known-class and vice versa. It is shown that both OSVM and 1-vs-Set SVM make imbalanced predictions by labelling too many known samples as unknown (90.91% and 95.33%), while our method has much better overall performances.

Method	overall	known	unknown	u → k	k → u
OSVM	58.12	9.09	99.89	0.11	90.91
1-vs-Set SVM	55.93	4.67	99.84	0.16	95.33
LSVM w/o outlier detection	69.19	63.41	74.12	25.88	17.80
LSVM w/ outlier detection (ours)	85.64	81.57	89.10	10.90	10.11

Table 2. The effect of pseudo-labeling process(%).

4. Effect of Number of Labeled Samples

In *Office* experiment, we set the number of labeled samples added from each labeled known categories to 3. We also analysis the impact of this number of labeled samples per class by conducting CDA with the number of labeled samples ranging from 1 to 9. The result on the three *Office* tasks are shown in Fig. 1.

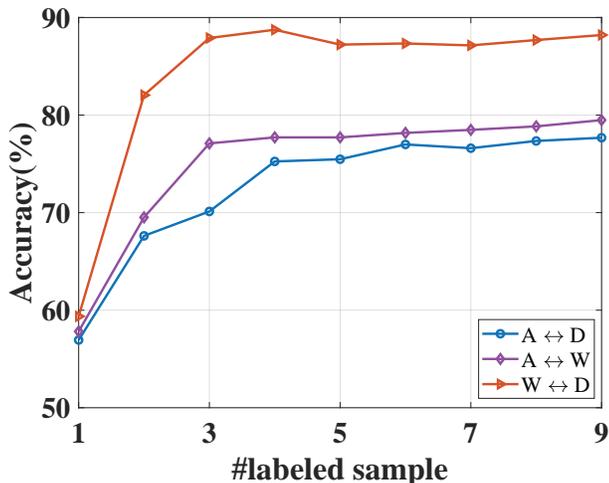


Figure 1. Impact of labeled sample number per class

The result showed that the performance of CDA gradually improves when more labeled samples are provided. This is expected because CDA use these labeled samples as initial inputs to assign pseudo labels for the unlabeled samples, and larger amount of labeled samples can produce more robust pseudo label prediction.

5. Effect of Overlapping Rate

In previous experiments, we set the overlapping rate of known label spaces between the two domains to 50% (e.g.,

Overlapping rate	0%	25%	50%	75%	100%
TCA	71.1	70.1	73.3	65.4	65.2
ATI-semi	73.8	72.5	73.4	73.8	74.8
MMDT	75.5	75.8	72.6	73.5	74.6
CDA	76.8	77.8	77.1	77.6	80.0

Table 3. Impact of the overlapping proportion(%).

class 1-10 for \mathcal{D}_A , 6-15 for \mathcal{D}_B). To compare the performance of our method under different overlapping rates, we have conducted experiments with overlapping rate varying from 0 to 1. We show the results of top competitors on the *Office* task A ↔ W in Table 3, and CDA performed the best under different circumstances.

6. Additional Compared Methods

Besides the methods compared in the paper, we have also compared with other domain adaptation methods. We first compare with 3 traditional methods: JDA [16], SA [5] and JGSA [26]. Then, 4 Deep Learning methods are compared as well, including: RevGrad [6], DAN [14], DCORAL [22] and DDC [24].

The full results of these methods as well as the methods included in the paper are shown in Table 4 and Table 5, for *Office* and DukeMTMC-reID respectively.

Methods	A ↔ W	A ↔ D	W ↔ D	AVG.
NA-avg	62.2 ± 3.19	61.0 ± 3.05	66.3 ± 1.72	63.12
NA	72.6 ± 2.04	69.4 ± 2.22	84.2 ± 3.89	75.40
TCA [18]	73.3 ± 1.67	70.6 ± 1.99	83.5 ± 3.96	75.80
GFK [7]	56.9 ± 2.89	55.7 ± 1.46	70.9 ± 4.85	61.17
CORAL [21]	69.9 ± 4.41	67.7 ± 3.13	83.7 ± 3.65	73.77
JAN [17]	63.8 ± 1.27	65.5 ± 0.76	74.7 ± 1.41	68.00
PADA [2]	60.3 ± 1.09	60.2 ± 0.98	70.9 ± 1.88	63.80
ATI [1]	70.4 ± 4.15	65.9 ± 1.80	81.7 ± 4.74	72.67
OpenBP [19]	62.6 ± 4.11	62.9 ± 1.71	67.9 ± 2.31	64.47
JDA [16]	72.6 ± 2.21	70.0 ± 2.03	83.2 ± 4.19	75.27
SA [5]	72.2 ± 2.07	69.9 ± 2.18	82.6 ± 3.87	74.90
JGSA [26]	62.3 ± 1.97	62.5 ± 1.05	67.2 ± 4.14	64.00
RevGrad [6]	57.1 ± 1.75	56.9 ± 1.60	58.0 ± 1.01	57.33
DDC [24]	55.1 ± 0.82	55.2 ± 0.94	55.6 ± 0.93	55.30
DAN [14]	69.4 ± 2.42	66.5 ± 4.13	83.6 ± 3.22	73.17
DCORAL [22]	68.6 ± 1.50	67.6 ± 4.02	83.5 ± 2.71	73.23
ATI-semi [1]	73.4 ± 2.28	72.0 ± 2.87	77.8 ± 3.40	74.72
MMDT [8]	72.6 ± 2.15	69.4 ± 2.19	84.4 ± 3.99	75.47
AMTL [12]	50.2 ± 1.45	48.8 ± 0.90	62.1 ± 2.17	53.70
CLMT [3]	50.3 ± 1.47	50.0 ± 0.77	61.7 ± 1.23	54.00
MRN [15]	62.2 ± 3.38	62.4 ± 2.44	77.4 ± 3.43	67.33
CDA	77.1 ± 1.35	75.2 ± 1.63	88.1 ± 2.45	80.13

Table 4. Comparing state-of-the-art methods on *Office*. The 1st/2nd best results are indicated in red/blue.

Method	1 ↔ 2		2 ↔ 3		3 ↔ 4		4 ↔ 5		5 ↔ 6		6 ↔ 7		7 ↔ 8		8 ↔ 1		AVG.	
	r=1	r=5																
NA-avg	43.6	55.3	29.6	38.2	50.1	54.8	69.2	95.2	38.2	54.4	33.5	42.4	22.0	40.3	76.7	95.4	45.4	59.5
NA	59.3	75.1	41.4	53.2	64.4	78.8	74.1	98.9	49.0	67.1	51.9	65.4	27.6	49.6	78.3	98.9	55.7	73.4
TCA [18]	58.6	74.7	41.0	53.7	64.3	79.0	75.3	98.9	48.6	67.3	51.5	65.3	27.2	49.8	78.3	98.9	55.6	73.5
GFK [7]	59.1	75.1	41.5	53.5	64.5	79.0	75.3	99.0	49.0	67.4	51.9	65.5	27.7	50.3	78.5	98.9	55.9	73.6
CORAL [21]	58.9	75.8	41.6	54.2	64.2	78.4	74.5	98.8	47.7	66.3	51.5	64.6	26.5	48.6	78.4	98.8	55.4	73.2
JAN [17]	24.0	43.7	34.4	64.4	21.4	38.6	75.5	90.2	30.2	60.0	27.7	52.4	44.6	72.1	81.7	92.5	42.4	64.2
PADA [2]	14.0	30.7	39.4	62.2	22.1	35.3	75.4	89.0	30.1	57.4	27.2	50.4	47.4	70.9	77.6	91.1	41.6	60.9
ATI [1]	58.6	74.0	40.6	49.5	65.4	79.2	36.1	78.0	47.9	63.6	52.4	65.6	24.6	42.7	28.2	84.4	44.2	67.1
OpenBP [19]	34.7	49.7	45.9	66.5	27.3	39.8	77.1	90.6	8.1	26.5	37.6	53.5	47.5	67.3	83.5	93.3	45.2	54.9
JDA [16]	58.9	73.4	39.6	48.6	66.9	79.5	43.7	84.0	45.2	59.1	49.7	60.7	22.8	36.0	35.9	84.2	45.3	65.7
SA [5]	59.3	75.1	41.4	53.2	64.4	78.8	74.1	98.9	49.0	67.1	51.9	65.4	25.7	49.7	78.3	99.0	55.5	73.4
JGSA [26]	45.7	61.7	35.1	48.0	62.7	74.7	16.9	88.7	40.3	53.5	43.2	55.6	16.5	31.3	7.2	36.5	33.5	56.2
RevGrad [6]	53.6	71.9	35.9	50.5	54.6	68.8	67.5	97.4	47.8	68.5	45.1	65.7	31.0	60.6	87.6	96.3	52.9	72.5
DAN [14]	53.2	71.0	35.0	47.6	54.9	68.9	77.8	97.2	45.5	66.3	45.6	65.2	31.3	59.9	82.6	97.2	53.2	71.7
DCORAL [22]	52.2	70.8	36.4	50.3	55.0	69.1	78.5	97.4	48.8	70.3	46.7	66.2	31.8	61.2	89.2	96.5	54.8	72.7
DDC [24]	53.1	72.4	37.5	52.7	55.4	71.3	74.2	97.7	46.4	66.7	45.2	63.8	30.6	58.3	92.3	95.6	54.3	72.3
MRN [15]	26.2	46.5	19.2	32.8	34.0	49.6	74.9	94.4	25.9	49.9	24.4	43.1	19.9	50.2	53.6	85.1	34.8	56.5
ATI-semi [1]	53.4	71.8	55.6	71.8	56.5	76.8	39.7	88.9	55.3	69.3	57.6	76.9	46.7	65.6	30.6	86.3	49.4	75.9
MMDT [8]	36.0	51.1	35.7	51.5	57.5	73.2	74.2	98.9	29.2	53.2	22.0	39.5	19.1	43.5	76.5	98.6	42.9	62.3
LMNN [25]	59.4	78.1	42.5	59.1	61.8	74.1	84.7	96.3	53.3	72.2	50.8	66.1	33.5	60.8	90.2	97.7	59.5	75.5
KISSME [11]	55.2	68.6	49.2	71.0	63.1	75.3	86.0	94.9	57.6	74.5	55.3	71.6	45.8	71.6	22.0	56.7	54.3	73.0
XQDA [13]	55.6	68.9	49.1	71.0	63.1	75.3	86.0	94.9	57.6	74.5	55.3	71.6	45.8	71.6	48.2	86.0	57.6	76.7
DLLR [10]	59.2	75.0	41.3	52.9	64.1	78.1	72.9	98.7	48.7	66.9	51.8	64.5	27.9	49.6	78.5	99.0	55.6	73.1
SPGAN [4]	41.6	59.9	31.3	47.5	44.7	58.6	69.6	97.7	41.4	61.0	43.3	59.6	22.8	46.3	74.1	98.0	48.7	69.6
CDA	64.6	80.4	62.4	88.1	67.9	84.9	76.0	98.9	61.5	82.1	59.6	78.1	67.2	83.8	85.6	98.5	68.1	86.9

Table 5. Comparing CMC accuracies with state-of-the-art methods on DukeMTMC-reID(%).

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