Curvature Diversity-Driven Deformation and Domain Alignment for Point Cloud

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Abstract

Unsupervised Domain Adaptation is crucial for point cloud learning due to geometric variations across different generation methods and sensors. To tackle this challenge, we propose **Curvature Diversity-Driven Nuclear-Norm Wasserstein Domain Alignment (CDND)**. We first introduce a Curvature Diversity-driven Deformation Reconstruction (CurvRec) task, enabling the model to extract salient features from semantically rich regions of a given point cloud. We then propose a theoretical framework for Deformation-based Nuclear-norm Wasserstein Discrepancy (D-NWD), extending the Nuclear-norm Wasserstein Discrepancy to original and deformed samples. Our theoretical analysis demonstrates that D-NWD is effective for any deformation method. Empirical experiment results show that our CDND achieves state-of-the-art performance by a noticeable margin over existing approaches.

1 Introduction

Adopting deep neural networks (DNNs) on point cloud representation learning has led to significant success in various applications, including robotics Maturana & Scherer (2015); Duan et al. (2021), autonomous vehicles Mahjourian et al. (2018); Cui et al. (2021), and scene understanding Zheng et al. (2013); Zhu et al. (2017). Point clouds captured under varying settings can exhibit significant differences, resulting in performance drops when DNNs are tested on data from settings different from the training set. In real-world scenarios, this discrepancy is known as the domain gap. The naive solution is to retrain the model using a new training dataset. However, obtaining labeled training data is both time-consuming and labor-intensive. Unsupervised domain adaptation (UDA) offers a solution to tackle these issues by utilizing knowledge transfer from source domains with annotated data to target domains with only unlabeled data. Although UDA is well studied for 2D planner data, e.g., images, UDA for 3D point clouds has not been explored extensively due to challenges such as the irregular, unstructured, and unordered nature of 3D point cloud data. Such irregularities exacerbate geometric variations between the source and target domains compared to the 2D planner data and make extending existing solutions nontrivial.

To address the above challenges, we propose Curvature Diversity-Driven Nuclear-Norm Wasserstein Domain Alignment (CDND). Our first contribution is to develop a deformation reconstruction method that leverages *curvature diversity* in different regions of a point cloud for domain alignment. We evaluate curvature diversity based on the entropy that captures the saliency of each region. This metric is then used to select regions for deformation and reconstruction. Unlike previous methods such as Achituve et al. (2021), which randomly select regions, our approach strategically selects them based on information content. Distinct from Zou et al. (2021), which classifies selected high-curvature regions into a fixed set, we focus on deforming and reconstructing regions with low curvature diversity. Our method avoids deforming semantically rich regions and enables the feature extractor to focus on extracting features from these regions for improved generalization.

Our second contribution is a theoretical framework for the Deformation-based Nuclear-norm Wasserstein Discrepancy (D-NWD). Unlike the conventional Nuclear-norm Wasserstein Discrepancy (NWD), D-NWD incorporates features from both original and deformed samples to align the source and target domains. By including features from deformed samples, D-NWD may create a diverse and robust feature space, enhancing

model generalization under domain shifts. Notably, prior work has not provided a theoretical analysis demonstrating that using the NWD to align features from both original and deformed (or augmented) samples can improve performance on the target domain dataset. Our analysis is the first to demonstrate the effectiveness of this strategy. Our theoretical framework shows that D-NWD effectively reduces the domain gap between source and target domains. Furthermore, our analysis illustrates that D-NWD is *generic* and adaptable to any deformation method, not just the one presented in this paper. Our theoretical analysis is non-trivial since incorporating features from deformed samples requires us to establish entirely new bounds and prove that the NWD remains effective in this expanded feature space spanning both original and deformed samples. Experiments on common benchmarks for both classification and segmentation show that our approach achieves SOTA performance compared to existing approaches.

2 Related Works

UDA for Point Clouds Despite extensive works on UDA for 2D planner image data Ganin & Lempitsky (2015); Mansour et al. (2008), only a limited number of studies Qin et al. (2019); Achituve et al. (2021); Shen et al. (2022); Zou et al. (2021); Liang et al. (2022); Chen et al. (2023); Katageri et al. (2024); Wei et al. (2024) address UDA in point clouds and non-planner data spaces as extending methods for 2D data for point clouds in non-trivial. Qin *et al.* Qin et al. (2019) introduce PointDAN that integrates both local and global domain alignment strategies. They also provide the PointDA benchmark for point cloud classification under the UDA setting. Achituve *et al.* Achituve et al. (2021) propose a domain alignment technique that involves reconstruction from deformation and incorporates PointMixup Chen et al. (2020). They also introduce the PointSegDA benchmark for point cloud segmentation under the UDA setting. Zou *et al.* Zou et al. (2021) utilize two geometry-inspired self-supervised classification tasks to learn domain-invariant feature. Shen *et al.* Shen et al. (2022) introduce a self-supervised method for learning geometry-aware implicit functions to handle domain-specific variations effectively. Liang *et al.* Liang et al. (2022) propose a masked local 3D structure prediction task that helps the model learn domain-invariant features by recovering masked parts of the point cloud. Katageri *et al.* Katageri et al. (2024) integrate contrastive learning with optimal transport to align features across domains, improving domain consistency and discrimination simultaneously. Chen *et al.* Chen et al. (2023) introduce a self-supervised boundary point prediction task to encourage the model to focus on geometrically informative boundary regions, which are often robust to domain shifts. Wei *et al.* Wei et al. (2024) develop a multi-scale part-based feature learning strategy to enhance both domain generalization and adaptation by capturing local and global part-level information. These methods show the growing interest in combining geometric cues with domain adaptation objectives. Our work differs by introducing a principled deformation strategy based on curvature diversity, a sophisticated self-supervised learning task, and a theoretical framework (D-NWD) that generalizes to any deformation method, leading to state-of-the-art performance.

Optimal Transport for Domain Adaptation The Wasserstein metric, known for encoding the natural geometry of probability measures within optimal transport theory, has been extensively studied for its application in domain adaptation due to its nice properties. Gautheron *et al.* Gautheron et al. (2019) propose Wasserstein Distance Guided Representation Learning to leverage the Wasserstein distance to enhance similarities between embedded features. Methods introduced in Lee et al. (2019) and Gabourie et al. (2019) use the sliced Wasserstein discrepancy instead of L_1 distance in Maximum Classifier Discrepancy Saito et al. (2018) to achieve a more geometrically meaningful intra-class divergence. Additionally, CGDM Du et al. (2021) introduces cross-domain gradient discrepancy to further mitigate domain differences. DeepJ-DOT Damodaran et al. (2018) utilizes a coupling matrix to map source samples to the target domain. Gautheron *et al.* Gautheron et al. (2019) propose a feature selection technique that addresses the domain shifts problem. Moreover, Xu *et al.* Xu et al. (2020) develop reliable weighted optimal transport, which uses spatial prototypical information and intra-domain structure to evaluate sample-level domain discrepancies, resulting in a better pairwise optimal transport plan. Finally, Fatras *et al.* Fatras et al. (2021) present an unbalanced optimal transport method combined with a mini-batch strategy to efficiently learn from large-scale datasets. In this work, we developed our D-NWD based on another previous method, NWD Chen et al. (2022).

3 Proposed Method

We begin by defining the unsupervised domain adaptation (UDA) problem, followed by an overview of our UDA approach, called Curvature Diversity-Driven Nuclear-Norm Wasserstein Domain Alignment (CDND), in Section 3.1. We then present the details of our main contributions: (1) a Curvature Diversity-based Deformation Reconstruction task, described in Sections 3.2 and 3.3, and (2) a theoretical framework for D-NWD, which extends NWD to incorporate both original and deformed samples, as detailed in Section 3.4 and 4. Note that our theoretical analysis is applicable to any deformation method used.

3.1 Problem Formulation

We consider a source domain with labeled samples and a target domain with unlabeled samples which has a different data distribution. Our goal is to develop a UDA method to train a model that accurately predicts labels for the target domain using both the source labeled dataset and the target unlabeled dataset. Let $\mathcal{S}$ represent the source domain, where X_s^i denotes the i-th batch of samples and y_s^i their corresponding labels. Similarly, let $\mathcal{T}$ represent the target domain, where X_t^i is the i-th batch of samples. The feature space induced by $\mathcal{S}$ and $\mathcal{T}$ is denoted by Ω_o. In addition, we introduce deformed domains $\mathcal{S}^d$ and $\mathcal{T}^d$, with their feature space Ω_d. We assume that Ω_o and Ω_d are disjoint subsets of $\mathbb{R}^n$, i.e., $\Omega_o \cap \Omega_d = \emptyset$, with $\Omega_o \cup \Omega_d \subseteq \mathbb{R}^n$. This assumption is generally valid in practice, as the probability of a deformed sample being exactly identical to a original one is negligible. A point cloud from the source domain is denoted as $x_s \in \mathbb{R}^{n\times 3}$ and from the target domain is $x_t \in \mathbb{R}^{n\times 3}$, where n is the number of points. The corresponding deformed point clouds are denoted by x_s^d and x_t^d.

The pipeline of our CDND is presented in Figure 1. Our model first uses a feature extractor E to obtain shape features from both source and target point clouds. To minimize domain gaps and ensure domain-invariant features, we: (1) use a curvature diversity-driven deformation reconstruction task via a reconstruction decoder h_{SSL} and (2) employ the D-NWD to align domains through a classifier C. The aligned features are then used for downstream tasks, *i.e.*, point cloud classification and segmentation. The model is trained using source-labeled and target-unlabeled data.

3.2 Curvature Diversity-Driven Deformation

To extract domain-invariant features from point clouds, three deformation strategies have been explored in the literature Achituve et al. (2021): volume-based, feature-based, and sample-based, according to the way of dividing point clouds into regions for deformation. Although these strategies use different techniques to select regions for deformation, they all divide a point cloud into regions based on some spatial locations or arrangements and uniformly randomly select regions to be deformed. However, uniformly random selection may not be optimal as regions within a point cloud vary in their semantic richness, *i.e.*, some regions contain more semantic information that is generalizable across the domains, while some regions may be domain-specific. These semantically rich regions are crucial for tasks such as classification, as they have more distinguishable characteristics. For instance, to differentiate a point cloud of a plant from that of a lamp, focusing on the leaves and flowers — which have richer semantic information — would be more effective than focusing on the flower pot, which is similar to the base of a lamp. Thus, deforming regions with richer semantic information causes the point cloud to lose semantic meaning, making it difficult for a classifier to classify it. Our idea is to identify shared semantic information to improve model generalization in the target domain. To encourage the feature extractor to prioritize regions with rich information, we propose deforming regions that are *less* semantically rich. Our deformation strategy helps the model learn to extract features from the most informative or salient regions of a point cloud.

To evaluate the richness of semantics, we propose using curvature diversity as a measurement. For the curvature diversity-driven deformation, we adopt the following steps. First, we use Farthest Point Sampling (FPS) Moenning & Dodgson (2003) to sample k center points as centers of k regions of a given point cloud. Then, for each center point, we use k-Nearest Neighbor (k-NN) to select m nearest points. As a result, each region is represented by a center point along with these m nearest points. Next, we select the N regions with the *smallest* curvature diversity to deform. To deform these selected regions, we replace all the points within

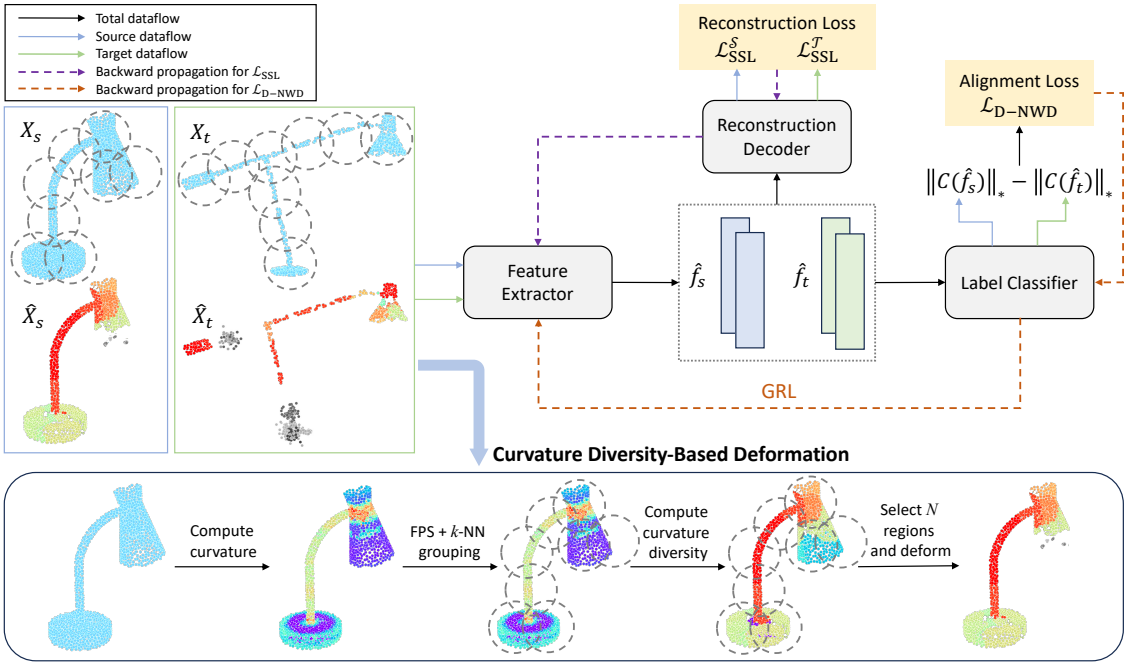

Figure 1: Pipeline of CDND. The inputs are the source batch X_s and target batch X_t. We first deform them into $\hat{X}_s$ and $\hat{X}_t$ using Curvature Diversity-Based Deformation. Next, X_s, X_t, $\hat{X}_s$, and $\hat{X}_t$ are sent into a feature extractor. The features of deformed samples are fed into a reconstruction decoder to reconstruct the deformed regions. For domain alignment, both original and deformed features are sent to D-NWD. Aside from the two losses shown in the figure, a cross-entropy loss is computed on X_s and $\hat{X}_s$ with labels. An NWD loss $\mathcal{L}_{\text{NWD}}^{\mathcal{T}}$ on X_t and $\hat{X}_t$ is also computed to ensure prediction consistency between the target original and deformed pairs.

these regions with new points. These new points are generated by sampling from a Gaussian distribution, where the mean is set to the average position of all the original points in that region, and the variance is set to 0.001. In Figure 1, $\hat{X}_s$ and $\hat{X}_t$ represent the deformed samples, and the points shown in grayscale are those drawn from the Gaussian distribution.

We compute point curvature using PCA Zou et al. (2021). Specifically, we first select a small neighborhood around each point and apply PCA to determine the principal directions and their eigenvalues. The curvature is then calculated as $c = |\lambda_{\min}| / \sum_{i=1}^{K} |\lambda_i|$, where $\lambda_{\min}$ is the smallest eigenvalue of the matrix, and K is the number of eigenvalues. Larger variation in curvature indicates a more intricate geometry and more significant shape changes within a region. The fourth lamp sample in the bottom row of Figure 1 illustrates this property: regions with warmer colors represent areas of higher curvature diversity. To measure the diversity or variation of curvature in a region, we propose to use *entropy of curvature*. Entropy effectively captures the variations in curvature values, allowing quantifying the richness of semantics within a region. Formally, we use the following to measure the curvature diversity:

$$
\begin{aligned}
c_{\min}^{j} &= \min_{c_i^j \in R^j} \{c_i^j\}_{i=1}^{N_{R^j}}, \quad c_{\max}^{j} = \max_{c_i^j \in R^j} \{c_i^j\}_{i=1}^{N_{R^j}}, \\
c_{i,\text{norm}}^{j} &= \frac{c_i^j - c_{\min}^j}{c_{\max}^j - c_{\min}^j + 10^{-10}}, \quad H(c_{\text{norm}}^j) = -\sum_{i=1}^{N_{R^j}} c_{i,\text{norm}}^j \cdot \log(c_{i,\text{norm}}^j + 10^{-10}),
\end{aligned}
\tag{1}
$$

where, c_i^j represents the curvatures of the i-th point in the j-th region of the point cloud which contains N_{R^j} points in total. To standardize these values, we first calculate $c_{\min}^j$ and $c_{\max}^j$, which are the minimum and maximum values of all curvatures within the j-th region, respectively. Using these values, we then normalize the curvature values to be in [0, 1], denoted as $\{c_{i,\text{norm}}^j\}$. Then, we calculate the curvature

diversity $H(c^j_{\text{norm}})$ by applying entropy.[1] The point-wise curvature values are computed only once before training. During training, we only compute the curvature diversity (entropy) for each region, based on these precomputed curvature values. This avoids repeated curvature estimation and keeps the runtime overhead minimal. Since curvature diversity calculation involves simple normalization and entropy over small regions, its cost is negligible compared to the main training process.

3.3 Deformation Reconstruction Loss

After deforming the selected regions, we obtain a deformed point cloud x^d from the original x. The deformed input x^d is processed by the feature extractor E to generate $E(x^d)$, which is then passed to a reconstruction decoder h_{SSL} to reconstruct x. The self-supervised loss $\mathcal{L}_{\text{SSL}}$ minimizes the distance between $h_{\text{SSL}}(E(x^d))$ and x. We use the Chamfer distance in $\mathcal{L}_{\text{SSL}}$, focusing on the original points in x within the deformed region R and their reconstructions from x^d. Formally, let $I \subset \{1, 2, \ldots\}$ represent the indices of the points in $x \cap R$, and we define $\mathcal{L}_{\text{SSL}}$ as follows:

$$\mathcal{L}_{\text{SSL}} = \sum_{(x, x^d, I)} \left(\{x_i\}_{i \in I}, \{h_{\text{SSL}}(E(x^d))_i\}_{i \in I} \right), \tag{2}$$

where x_i is the i-th point in the point cloud x and D denotes the Chamfer distance which is defined as:

$$D(R_1, R_2) = \sum_{a \in R_1} \min_{b \in R_2} \|a - b\|_2^2 + \sum_{b \in R_2} \min_{a \in R_1} \|b - a\|_2^2, \tag{3}$$

where $D(R_1, R_2)$ measures the discrepancy between point cloud regions $R_1, R_2 \subset \mathbb{R}^3$. Note that when we reconstruct only the deformed regions, we can also reduce computational burden and improve time efficiency.

3.4 Domain Alignment via D-NWD

We begin with introducing the 1-Wasserstein distance.

Definition 1 (1-Wasserstein distance). Adler & Lunz (2018) *The 1-Wasserstein distance quantifies the minimal cost of transporting mass between two probability measures defined on the same metric space. Let μ and ν be two probability measures over a metric space (Ω, d), where $d(x, y)$ denotes the distance between points x and y in Ω. Then the 1-Wasserstein distance is defined as:*

$$W_1(\mu, \nu) = \inf_{\gamma \in \Gamma(\mu, \nu)} \int_{\Omega \times \Omega} d(x, y) \, d\gamma(x, y), \tag{4}$$

where $\Gamma(\mu, \nu)$ is the set of all couplings with marginals μ and ν; that is,

$$\int_\Omega \gamma(x, y) \, dy = \mu(x), \quad \int_\Omega \gamma(x, y) \, dx = \nu(y).$$

The Kantorovich–Rubinstein duality further states that the 1-Wasserstein distance can be expressed in a dual form as:

$$W_1(\mu, \nu) = \sup_{\|h\|_L \le K_L} \mathbb{E}_{x \sim \mu}[h(x)] - \mathbb{E}_{x \sim \nu}[h(x)], \tag{5}$$

where the supremum is taken over all functions $h : \Omega \to \mathbb{R}$ with Lipschitz constant at most K_L, i.e.,

$$\|h\|_L := \sup_{x \neq y} \frac{|h(x) - h(y)|}{d(x, y)} \le K_L. \tag{6}$$

The Nuclear-norm Wasserstein Discrepancy (NWD) Chen et al. (2022) belongs to the family of 1-Wasserstein distances, with sophisticatedly chosen h's. Below, we present the form of h in NWD. Consider a prediction

[1] Rigorously speaking, $\{c^j_{i,\text{norm}}\}$ is an un-normalized distribution without being divided by a partition function or normalization constant, but it does not affect our claim of curvature diversity.

matrix $P \in \mathbb{R}^{b \times M}$ predicted by the classifier C, where b is the number of samples in a batch and M is the number of classes. The non-negative self-correlation matrix $Z \in \mathbb{R}^{M \times M}$ is computed as $Z = P^T P$. The intra-class correlation I_a is defined as the sum of the main diagonal elements of Z, and the inter-class correlation I_e is the sum of all the off-diagonal elements of Z:

$$I_a = \sum_{i=1}^{M} Z_{ii}, \quad I_e = \sum_{i \neq j}^{M} Z_{ij}.$$

In the source domain, I_a is large, and I_e is relatively small because most samples are correctly classified. Conversely, in the target domain, I_a is small, and I_e is relatively large due to the large error from the lack of supervised training on the target domain. Hence, $I_a - I_e$ can represent the discrepancy between the two domains, as $I_a - I_e$ is large for the source domain but small for the target domain. Note that $I_a = \|P\|_F^2$ can be represented as the squared Frobenius norm of P, and thus $I_a - I_e = 2\|P\|_F^2 - b$. [2] We can rewrite $P_s = C(F_s)$ and $P_t = C(F_t)$, where F_s and F_t are feature representation batches from the source and target domains, respectively. We find $\|C\|_F$ gives high scores to the source domain and low scores to the target domain. Thus, we can set h in Eq. 5 to be $\|C\|_F$ and represent the domain discrepancy as:

$$W_F(\nu_s, \nu_t) = \sup_{\|\|C\|_F\|_L \leq K_L} \mathbb{E}_{F_s \sim \nu_s}[\|C(F_s)\|_F] - \mathbb{E}_{F_t \sim \nu_t}[\|C(F_t)\|_F],$$

where ν_s is the probability measure for features of samples in $\mathcal{S}$ and ν_t is the probability measure for features of samples in $\mathcal{T}$. To enhance prediction diversity, the Frobenius norm can be replaced with the nuclear norm, which maximizes the rank of P while still being bounded by the Frobenius norm Chen et al. (2022). Thus, the domain discrepancy can be rewritten as:

$$W_N(\nu_s, \nu_t) = \sup_{\|\|C\|_*\|_L \leq K_L} \mathbb{E}_{F_s \sim \nu_s}[\|C(F_s)\|_*] - \mathbb{E}_{F_t \sim \nu_t}[\|C(f_t)\|_*]. \tag{7}$$

The Eq. 7 is the formal definition of NWD, where C denotes the classifier, and $\|\cdot\|_*$ represents the nuclear norm. In the NWD paper, empirically, NWD can be approximate by $\mathcal{L}_{\text{NWD}}$:

$$\mathcal{L}_{\text{NWD}} = \frac{1}{B_s} \sum_{i=1}^{B_s} \|C(F_s^i)\|_* - \frac{1}{B_t} \sum_{i=1}^{B_t} \|C(F_t^i)\|_*$$

where $F_s^i \sim \nu_s$ are the features for the i-th source batch and $F_t^i \sim \nu_t$ are the features for the i-th target batch. B_s is the number of training batches in the source domain and B_t is the number of training batches in the target domain.

$$\min_E \max_C \mathcal{L}_{\text{NWD}} \tag{8}$$

Then, the domain alignment is achieved through a min-max game presented in Eq. 8.

Now, we introduce our second component, Deformation-based Nuclear-norm Wasserstein Discrepancy (D-NWD). From the previous sections, the curvature diversity-driven deformation reconstruction helps reduce the domain gap between the source and target domains. To further complete classification or segmentation tasks in the presence of the domain gap, we propose D-NWD to align domains, as inspired by NWD Chen et al. (2022). Our D-NWD objective is defined as:

$$W_N(\nu_{s \cup s^d}, \nu_{t \cup t^d}) = \sup_{\|\|C\|_*\|_L \leq K_L} \mathbb{E}_{\hat{F}_s \sim \nu_{s \cup s^d}}[\|C(\hat{F}_s)\|_*] - \mathbb{E}_{\hat{F}_t \sim \nu_{t \cup t^d}}[\|C(\hat{F}_t)\|_*], \tag{9}$$

Here, $\nu_{s \cup s^d}$ and $\nu_{t \cup t^d}$ are probability measures defined over $\Omega_o \cup \Omega_d$, for the features from samples in the original and deformed source and target domains. We align the probability measure of features from the original and deformed samples in the source domain with that of the target domain. Our motivation is that taking features from deformed samples into account would provide a richer, more robust feature space, reduce overfitting, and increase the model's adaptability to variations inherent in data. This technique differs from

[2] We have $\sum_{j=1}^{M} Z_{i,j} = 1$, $\forall i \in \{1, \cdots, b\}$ and $j \in \{1, \cdots, M\}$, and thus $I_a + I_e = b$ Chen et al. (2022).

using NWD, which aligns ν_s and ν_t defined over Ω_o, the probability measures for the features from samples in the original source and target domains. Empirically, our objective in Eq. 9 be approximated by $\mathcal{L}_{\text{D-NWD}}$:

$$\mathcal{L}_{\text{D-NWD}} = \frac{1}{B_s}\sum_{i=1}^{B_s} \|C(\hat{F}_s^i)\|_* - \frac{1}{B_t}\sum_{i=1}^{B_t} \|C(\hat{F}_t^i)\|_*, \tag{10}$$

$\hat{F}_s^i \sim \nu_{s\cup s^d}$ represents the features for the i-th source batch and $\hat{F}_t^i \sim \nu_{t\cup t^d}$ represents the features for the i-th target batch. In practice, we obtain the original and deformed samples by first sampling from the original domain and then generating the corresponding deformed versions. The alignment is then performed through a min-max game, described in the following:

$$\min_E \max_C \mathcal{L}_{\text{D-NWD}}. \tag{11}$$

To avoid alternating updates, we employ a Gradient Reverse Layer Ganin et al. (2016), following the approach in Chen et al. (2022), to make the learned features discriminative and domain-agnostic.

3.5 Overall Loss

In addition to deformation and domain alignment loss defined in Eq. 2 and Eq. 11, we use a cross-entropy loss $\mathcal{L}_{\text{CLS}}$ on both the original and the deformed source domain samples for supervised training of the classifier:

$$\mathcal{L}_{\text{CLS}} = \frac{1}{B_s}\sum_{i=1}^{B_s} \mathcal{L}_{\text{CE}}(C(\hat{F}_s^i), Y_s^i), \tag{12}$$

where Y_s^i are labels for batch $\hat{F}_s^i$. Since we have no access to the ground-truth labels for the target domain data, it is impossible to use the supervised cross-entropy loss as in Eq. 12 on samples from $\mathcal{T}$ and $\widetilde{\mathcal{T}}$. One straightforward alternative is to adopt pseudo-labels as in Fan et al. (2022); Liang et al. (2022); Zou et al. (2021); Shen et al. (2022). However, this strategy has the risk that the classifier might mistakenly predict target samples as the major classes of the source domain. Instead, we use NWD to ensure consistency in predictions between $\mathcal{T}$ and $\mathcal{T}^d$. Thus, we define a target domain loss $\mathcal{L}_{\text{NWD}}^{\mathcal{T}}$ as:

$$\mathcal{L}_{\text{NWD}}^{\mathcal{T}} = \frac{1}{B_t}\sum_{i=1}^{B_t} \|C(F_t^i)\|_* - \frac{1}{B_t}\sum_{i=1}^{B_t} \|C(F_{t^d}^i)\|_*, \tag{13}$$

where $F_{t^d}^i \sim \nu_{t^d}$ denotes the deformed target domain batch and $F_t^i \sim \nu_t$ denotes the original target domain batch. Combining Eq. 2, Eq. 12, Eq. 11 and Eq. 13 together, our overall objective loss is:

$$\begin{aligned} &\min_{E,h_{\text{SSL}},C} \alpha\mathcal{L}_{\text{CLS}} + \gamma\mathcal{L}_{\text{SSL}}, \\ &\min_E \max_C \beta_1\mathcal{L}_{\text{D-NWD}} + \beta_2\mathcal{L}_{\text{NWD}}^{\mathcal{T}}, \end{aligned} \tag{14}$$

where $\alpha, \gamma, \beta_1, \beta_2$ can be tuned using the target domain validation set and setting details can be found in Appendix A.2. Note that $\mathcal{L}_{\text{D-NWD}}$ and $\mathcal{L}_{\text{NWD}}^{\mathcal{T}}$ serve distinct and complementary purposes in Equation 14. $\mathcal{L}_{\text{D-NWD}}$ is specifically used to minimize the domain gap between source and target features, while $\mathcal{L}_{\text{NWD}}^{\mathcal{T}}$ is used to improve prediction consistency between original and deformed target distributions, functioning similarly to a cross-entropy loss between pseudo-labels and predicted labels on target samples. We use NWD over pseudo-labeling since the latter can reinforce incorrect predictions when the classifier is biased toward majority source domain classes.

4 Theoretical Analysis

In this section, we present our theoretical contribution for D-NWD, demonstrating the effectiveness of extending NWD to both original and deformed samples. We provide new bounds in Theorems 1 and 2

along with their non-trivial proofs. Following Ben-David et al. (2006) and Chen et al. (2022), we perform our analysis in a binary classification scenario, which can be easily adapted to multi-class classification through reduction techniques such as one-vs-all Rifkin & Klautau (2004) or one-vs-one Allwein et al. (2000) approaches. Consider $\{C : \mathbb{R}^n \to [0,1]\}$ as a set of source classifiers within the hypothesis space $\mathcal{H}$. Let ν_s defined on Ω_o be the probability measure of original source domain and ν_{s_d} defined on Ω_d be the probability measure of deformed source domain. We define ν_t and ν_t^d in a similar way. The risk or error of classifier C on the original source domain is defined as $\varepsilon_s(C) = \mathbb{E}_{f_s \sim \nu_s}[|C(f_s) - y_s|]$, where y_s is the label associated with the feature f_s. We then define $\varepsilon_{s \cup s^d}(C) = \mathbb{E}_{\hat{f}_s \sim \nu_{s \cup s^d}}[|C(\hat{f}_s) - \hat{y}_s|]$, where $\hat{y}_s$ is the label associated with $\hat{f}_s$. Similarly, we define $\varepsilon_t(C), \varepsilon_{t \cup t^d}(C)$ as the errors on the target domain. The optimal classifier is defined as $C^* = \arg\min_C \varepsilon_{s \cup s^d}(C) + \varepsilon_t(C)$ which minimizes the combined error across $\nu_{s \cup s^d}$ and ν_t. Our Theorem 1 demonstrates that the expected target error $\varepsilon_t(C)$ can be bounded by the D-NWD on $\nu_{s \cup s^d}$ and $\nu_{t \cup t^d}$, $W_N(\nu_{s \cup s^d}, \nu_{t \cup t^d})$. Building on Theorem 1, we derive Theorem 2. Theorem 2 establishes that $\varepsilon_t(C)$ can be bounded by D-NWD on empirical probability measures $\hat{\nu}_{s \cup s^d}$ and $\hat{\nu}_{t \cup t^d}$, $W_N(\hat{\nu}_{s \cup s^d}, \hat{\nu}_{t \cup t^d})$. **All proofs can be found in Appendix A.1.**

Theorem 1. *Let $(\Omega_o, \mathcal{F}_o, \nu_s)$, $(\Omega_d, \mathcal{F}_d, \nu_{s^d})$, $(\Omega_o, \mathcal{F}_o, \nu_t)$, and $(\Omega_d, \mathcal{F}_d, \nu_{t^d})$ be four probability spaces, where Ω_o and Ω_d are disjoint and $\Omega_o \cup \Omega_d \subseteq \mathbb{R}^n$. With the results of Lemma 1, let $(\Omega_o \cup \Omega_d, \mathcal{F}_u, \nu_{s \cup s^d})$ and $(\Omega_o \cup \Omega_d, \mathcal{F}_u, \nu_{t \cup t^d})$ be two probability spaces with probability measures defined as $\nu_{s \cup s^d} = 1/2\nu_s + 1/2\nu_{s^d}$ and $\nu_{t \cup t^d} = 1/2\nu_t + 1/2\nu_{t^d}$. Specifically, when sampling from $\nu_{t \cup t^d}$, there is an equal probability of $1/2$ to sample from ν_t or ν_{t^d}. Similarly, sampling from $\nu_{s \cup s^d}$ gives an equal probability of $1/2$ to draw from ν_s or ν_{s^d}. Consider a classifier $C \in \mathcal{H}_1$ and an ideal classifier $C^* = \arg\min_C \varepsilon_{s \cup s^d}(C) + \varepsilon_t(C)$ satisfying the K_L-Lipschitz constraint, where $\mathcal{H}_1$ is a subspace of the hypothesis space $\mathcal{H}$. For every classifier C in $\mathcal{H}_1$, the following inequality holds:*

$$\varepsilon_t(C) \leq 2\varepsilon_{s \cup s^d}(C) + 4K_L \cdot W_N(\nu_{s \cup s^d}, \nu_{t \cup t^d}) + \eta^*, \tag{15}$$

where $\eta^* = 2\varepsilon_{s \cup s^d}(C^*) + \varepsilon_t(C^*)$ is the ideal combined error and is a sufficiently small constant.

Theorem 2. *Under the assumption of Theorem 1, Ω_o and Ω_d are disjoint and $\Omega_o \cup \Omega_d \subseteq \mathbb{R}^n$. Let $(\Omega_o \cup \Omega_d, \mathcal{F}_u, \nu_{s \cup s^d})$ and $(\Omega_o \cup \Omega_d, \mathcal{F}_u, \nu_{t \cup t^d})$ be two probability spaces with $\nu_{s \cup s^d} = 1/2\nu_s + 1/2\nu_{s^d}$ and $\nu_{t \cup t^d} = 1/2\nu_t + 1/2\nu_{t^d}$, where $\nu_s, \nu_{s^d}, \nu_t, \nu_{t^d}$ each has a square-exponential moment. From Lemma 3 and 4, $\nu_{s \cup s^d}$ satsifies $T_1(\eta_s)$ for some η_s and $\nu_{t \cup t^d}$ satsifies $T_1(\eta_t)$ for some η_t. Let $\{\hat{f}_s^i\}_{i=1}^{N_s}$ and $\{\hat{f}_t^i\}_{i=1}^{N_t}$ be two sample sets of size N_s and N_t drawn i.i.d from $\nu_{s \cup s^d}$ and $\nu_{t \cup t^d}$, respectively. $\hat{\nu}_{s \cup s^d} = \frac{1}{N_s}\sum_{i=1}^{N_s} \delta_{\hat{f}_s^i}$ and $\hat{\nu}_{t \cup t^d} = \frac{1}{N_t}\sum_{i=1}^{N_t} \delta_{\hat{f}_t^i}$ are associated empirical probability measures. Then, for any $n' > n$ and $\eta' < \min(\eta_s, \eta_t)$, there exists a constant N_0 depending on n' such that for any $\delta > 0$ and $\min(N_s, N_t) \geq N_0 \max(\delta^{-(n'+2)}, 1)$, with probability at least $1 - \delta$, the following holds for all C:*

$$\varepsilon_t(C) \leq 2\varepsilon_{s \cup s^d}(C) + 4K_L \cdot W_N(\hat{\nu}_{s \cup s^d}, \hat{\nu}_{t \cup t^d}) + \eta^* + 4K_L \cdot \sqrt{\frac{2}{\eta'} \log \frac{1}{\delta}} \left(\sqrt{\frac{1}{N_s}} + \sqrt{\frac{1}{N_t}} \right), \tag{16}$$

where $\eta^* = 2\varepsilon_{s \cup s^d}(C^*) + \varepsilon_t(C^*)$ is the ideal combined error and is a sufficiently small constant.

Equation 16 justifies why our method can be effective empirically. Specifically, η^* are sufficiently small constants for relevant domains with consistent labels because it is the error corresponding to the ideal classifier C^*. The term $\sqrt{\frac{2}{\eta'} \log \frac{1}{\delta}} \left(\sqrt{\frac{1}{N_s}} + \sqrt{\frac{1}{N_t}} \right)$ is also a small constant when training dataset sizes, N_s and N_t, are large. $\varepsilon_{s \cup s^d}(C)$ is minimized by a supervised classification loss, since source domain samples have labels. Therefore, the primary objective is to minimize $W_N(\hat{\nu}_{s \cup s^d}, \hat{\nu}_{t \cup t^d})$, our D-NWD on empirical measures $\hat{\nu}_{s \cup s^d}$ and $\hat{\nu}_{t \cup t^d}$. Hence, minimizing $W_N(\hat{\nu}_{s \cup s^d}, \hat{\nu}_{t \cup t^d})$ can reduce error on the original target domain $\mathcal{T}$ and improve the model's performance on the target domain. It is important to note that our theoretical analysis is not intended to prove that D-NWD is superior to NWD. A direct comparison between NWD and D-NWD is not feasible, as they apply to different probability measures: $\hat{\nu}_{s \cup s^d}, \hat{\nu}_{t \cup t^d}$ for D-NWD and $\hat{\nu}_s, \hat{\nu}_t$ for NWD. Our theoretical contribution lies in showing that, regardless of the deformation method used, optimizing D-NWD on $\hat{\nu}_{s \cup s^d}$ and $\hat{\nu}_{t \cup t^d}$ can effectively reduce error on samples from $\mathcal{T}$. In other words, D-NWD mitigates the negative effects of domain gaps and enhances performance on the original target domain $\mathcal{T}$. Note that

We use equal mixture weights—specifically, a 1:1 ratio—in the probability measure (e.g., $\frac{1}{2}\nu_s + \frac{1}{2}\nu_s^d$) to simplify the implementation. However, this assumption is not essential for the validity of our conclusions. The theoretical results can be generalized to arbitrary sampling ratios by adjusting the mixture weights in the probability measures accordingly. In such cases, the only change would be in the constant factors reflecting the updated contribution from each component in the mixture. The core theoretical framework and its implications remain intact. In Appendix A.3, we present the experiment results with different ratios.

5 Experiments

We evaluate our method on the **PointDA-10** Qin et al. (2019) dataset, a domain adaptation dataset for point cloud classification, and on **PointSegDA** Achituve et al. (2021), a dataset for point cloud segmentation. For the PointDA-10 dataset, we compare our approach against the state-of-the-art methods for point cloud domain adaptation, including **DANN** Ganin et al. (2016), **PointDAN** Qin et al. (2019), **RS** Sauder & Sievers (2019), **DefRec+PCM** Achituve et al. (2021), **GAST** Zou et al. (2021), **ImplicitPCDA** Shen et al. (2022), and the most recent method with publicly available code, **PCFEA** Wang & el al (2025). Note that many recent methods do not provide code for reproducibility, which limits their inclusion. For the **PointSegDA** dataset, we compare our method with **RS**, **DefRec+PCM**, **GAST**, **ImplicitPCDA**, and **Adapt-SegMap** Tsai et al. (2018). Unfortunately, no suitable methods from the past 1–2 years with available code are applicable to this dataset. For both datasets, we also evaluate two upper bounds: **Supervised-T**, which involves training exclusively on labeled target samples, and **Supervised**, which uses both labeled source and target samples. Additionally, we assess a lower bound, **Unsupervised**, which utilizes only labeled source samples.

Additionally, for the PointDA-10 dataset, we incorporate Self-Paced Self-Training (SPST) into GAST, ImplicitPCDA, and our method, as SPST is originally included in both GAST and ImplicitPCDA. We exclude SPST for the PointSegDA dataset. SPST typically relies on ranking training samples by difficulty and gradually incorporating harder examples into training. The training samples for point cloud segmentation tasks are points in point clouds. However, mIoU is a global metric that evaluates performance across an entire point cloud, making it challenging to assign difficulty scores to individual points in a point cloud. The mechanism of SPST mismatches the per-point cloud, rather than per-point, evaluation criterion of mIoU. **Hyperparameter settings and implementation details can be found in Appendix A.2.**

5.1 Datasets

We use **PointDA-10** and **PointSegDA** datasets in our experiments. **PointDA-10** consists of three domains: ShapeNet-10 Chang et al. (2015), ModelNet-10 Wu et al. (2015), and ScanNet-10 Dai et al. (2017), each sharing ten distinct classes. **PointSegDA** consists of four domains: ADOBE, FAUST, MIT, and SCAPE. These domains share eight distinct classes of human body parts but vary in point distribution, pose, and scanned humans.

5.2 Training Scheme

We use DGCNN as the feature extractor Achituve et al. (2021) for fair comparison. We repeat our experiments three times using distinct random seeds for initialization and report the average accuracy and standard deviation. To ensure a fair comparison, we maintain the same seed for data shuffling and use the Adam optimizer Kingma & Ba (2014) for optimization.

5.3 Comparative Results

Results on PointDA. The results are presented in Table 1. We use S^+ to represent the ScanNet dataset, M to represent ModelNet, and S to represent the ShapeNet dataset. MS, SM, etc. are the abbreviated source and target domain pairs. The CDND model shows significant improvement over the other approaches on the PointDA-10 dataset with the highest average accuracy of 70.3%, outperforming all other models. CDND delivers state-of-the-art performance on five out of six tasks. It excels in tasks with a large domain

Models	**MS**	**MS$^+$**	**SM**	**SS$^+$**	**S$^+$M**	**S$^+$S**	**Avg**
Supervised-T	$93.9_{\pm 0.2}$	$78.4_{\pm 0.6}$	$96.2_{\pm 0.1}$	$78.4_{\pm 0.6}$	$96.2_{\pm 0.4}$	$93.9_{\pm 0.2}$	89.5
Unsupervised	$83.3_{\pm 0.7}$	$43.8_{\pm 2.3}$	$75.5_{\pm 1.8}$	$42.5_{\pm 1.4}$	$63.8_{\pm 3.9}$	$64.2_{\pm 0.8}$	62.2
DANN Ganin et al. (2016)	$75.3_{\pm 0.6}$	$41.5_{\pm 0.2}$	$62.5_{\pm 1.4}$	$46.1_{\pm 2.8}$	$53.3_{\pm 1.2}$	$63.2_{\pm 1.2}$	57.0
PointDAN Qin et al. (2019)	$82.5_{\pm 0.8}$	$47.7_{\pm 1.0}$	$77.0_{\pm 0.3}$	$48.5_{\pm 2.1}$	$55.6_{\pm 0.6}$	$67.2_{\pm 2.7}$	63.1
RS Sauder & Sievers (2019)	$81.5_{\pm 1.2}$	$35.2_{\pm 5.9}$	$71.9_{\pm 1.4}$	$39.8_{\pm 0.7}$	$61.0_{\pm 3.3}$	$63.6_{\pm 3.4}$	58.8
DefRec+PCM Achituve et al. (2021)	$81.7_{\pm 0.6}$	$51.8_{\pm 0.3}$	$78.6_{\pm 0.7}$	$54.5_{\pm 0.3}$	$73.7_{\pm 1.6}$	$71.1_{\pm 1.4}$	68.6
PCFEA Wang & el al (2025)	$84.4_{\pm 1.1}$	$47.9_{\pm 2.9}$	$71.8_{\pm 1.9}$	$47.1_{\pm 0.7}$	$71.8_{\pm 7.4}$	$68.1_{\pm 2.1}$	65.2
GAST Zou et al. (2021)	$82.3_{\pm 0.6}$	$53.0_{\pm 1.1}$	$72.6_{\pm 1.9}$	$47.6_{\pm 1.5}$	$64.6_{\pm 1.5}$	$66.8_{\pm 0.6}$	64.5
GAST+SPST	$84.5_{\pm 0.5}$	$54.1_{\pm 1.8}$	$80.1_{\pm 4.6}$	$46.7_{\pm 0.6}$	$81.5_{\pm 1.7}$	$66.7_{\pm 1.1}$	68.9
ImplicitPCDA Shen et al. (2022)	$79.5_{\pm 0.4}$	$41.7_{\pm 1.3}$	$72.9_{\pm 1.0}$	$47.5_{\pm 2.9}$	$67.6_{\pm 5.2}$	$66.4_{\pm 0.9}$	62.6
ImplicitPCDA+SPST	$81.3_{\pm 2.2}$	$33.2_{\pm 13.4}$	$73.2_{\pm 3.4}$	$38.0_{\pm 4.6}$	$66.9_{\pm 7.7}$	$75.0_{\pm 2.7}$	61.3
CDND	$84.1_{\pm 0.3}$	$\mathbf{58.7}_{\pm 0.8}$	$76.2_{\pm 0.0}$	$\mathbf{55.7}_{\pm 1.0}$	$75.1_{\pm 1.5}$	$72.0_{\pm 1.9}$	70.3
CDND+SPST	$\mathbf{85.4}_{\pm 1.1}$	$57.6_{\pm 1.3}$	$\mathbf{85.0}_{\pm 2.2}$	$54.5_{\pm 1.1}$	$\mathbf{82.6}_{\pm 0.7}$	$\mathbf{74.6}_{\pm 4.4}$	**73.3**

Table 1: Performance results (accuracy) on PointDA-10 dataset.

Models	**MS**	**MS$^+$**	**SM**	**SS$^+$**	**S$^+$M**	**S$^+$S**	**Avg**
NWD	$83.3_{\pm 0.7}$	$46.7_{\pm 1.7}$	$75.5_{\pm 1.8}$	$48.9_{\pm 2.5}$	$63.8_{\pm 3.9}$	$66.7_{\pm 1.9}$	64.2
DefRec	$83.4_{\pm 0.5}$	$46.9_{\pm 2.3}$	$74.5_{\pm 0.9}$	$46.3_{\pm 0.6}$	$67.7_{\pm 2.3}$	$64.0_{\pm 0.8}$	64.0
DefRec+NWD	$83.4_{\pm 0.5}$	$51.2_{\pm 3.0}$	$74.5_{\pm 0.9}$	$53.7_{\pm 3.8}$	$67.7_{\pm 2.3}$	$68.5_{\pm 2.4}$	66.5
DefRec+D-NWD	$83.4_{\pm 0.5}$	$53.1_{\pm 2.3}$	$74.5_{\pm 0.9}$	$54.6_{\pm 1.0}$	$67.7_{\pm 2.3}$	$67.4_{\pm 0.1}$	66.8
CurvRec(S)-High	$83.8_{\pm 0.9}$	$52.0_{\pm 1.4}$	$\mathbf{78.0}_{\pm 1.0}$	$45.9_{\pm 3.8}$	$72.5_{\pm 1.4}$	$66.7_{\pm 1.1}$	66.5
CurvRec(S)-Low	$83.1_{\pm 0.9}$	$53.0_{\pm 1.9}$	$74.9_{\pm 0.8}$	$44.7_{\pm 1.2}$	$74.8_{\pm 0.9}$	$65.9_{\pm 0.2}$	66.1
CurvRec(En)-High	$82.9_{\pm 1.5}$	$52.1_{\pm 0.4}$	$77.0_{\pm 0.3}$	$46.7_{\pm 1.0}$	$70.9_{\pm 0.6}$	$65.8_{\pm 0.4}$	65.9
CurvRec	$\mathbf{84.1}_{\pm 0.3}$	$52.2_{\pm 1.3}$	$76.2_{\pm 0.0}$	$50.1_{\pm 0.3}$	$\mathbf{75.1}_{\pm 1.5}$	$66.4_{\pm 1.5}$	67.4
CurvRec+PCM	$83.0_{\pm 0.5}$	$53.7_{\pm 1.0}$	$74.0_{\pm 0.6}$	$54.8_{\pm 1.1}$	$73.8_{\pm 1.1}$	$\mathbf{76.8}_{\pm 0.9}$	69.4
CurvRec+NWD	$\mathbf{84.1}_{\pm 0.2}$	$54.3_{\pm 2.2}$	$76.2_{\pm 0.0}$	$52.7_{\pm 2.1}$	$\mathbf{75.1}_{\pm 1.5}$	$70.6_{\pm 2.2}$	68.8
CDND, $\beta_2 = 0$	$84.1_{\pm 0.3}$	$57.4_{\pm 1.2}$	$76.2_{\pm 0.0}$	$\mathbf{55.7}_{\pm 0.2}$	$75.1_{\pm 1.5}$	$69.9_{\pm 1.7}$	69.7
CDND (CurvRec+D-NWD)	$\mathbf{84.1}_{\pm 0.3}$	$\mathbf{58.7}_{\pm 0.8}$	$76.2_{\pm 0.0}$	$\mathbf{55.7}_{\pm 1.0}$	$\mathbf{75.1}_{\pm 1.5}$	$72.0_{\pm 1.9}$	**70.3**

Table 2: Ablation study results (accuracy) on PointDA-10 dataset.

gap, such as MS$^+$, S$^+$M, SS$^+$, and S$^+$S. In these tasks, one domain is a synthetic dataset, and the other domain is a real-world dataset. This shows its proficiency in handling complex transformations. Especially, CDND scores 58.7% on MS$^+$, outperforming the second-best method by approximately 6%. Additionally, CDND maintains competitive accuracy in tasks with a small domain gap, such as SM and MS, with scores of 84.1% and 76.2%, respectively. With SPST, the performance is further improved, as CDND+SPST achieves 73.3%, outperforming GAST+SPST by 4.4% and ImplicitPCDA+SPST by 12%. Note that plain CDND also outperforms both GAST+SPST and ImplicitPCDA+SPST on average. The performance of CDND across various tasks highlights its ability to adapt to diverse domain challenges, making it a promising choice for point cloud classification in the UDA setting.

Results on PointSegDA. The results are presented in Table 3 We use A to represent the ADOBE dataset, F to represent the FAUST dataset, M to represent the MIT dataset, and S to represent the SCAPE dataset. AF, FA, etc. are the abbreviated source and target domain pairs. On the PointSegDA benchmark, CDND achieves the highest average score of 59.9%, which surpasses the second-best method, RS, by a margin of 2.0%, which is significant in terms of mIoU on the segmentation task. Its superior performance is particularly evident in MA and SA tasks; in the MA task, CDND achieves a mIoU of 68.6, outperforming RS by 9%. Similarly, in the SA task, CDND secures a mIoU of 77.5, which is around 7% higher than RS. These results showcase its adaptability and learning capability. Additionally, in the FA task, CDND achieves a score of 81.5, even slightly surpassing the supervised baseline. In other tasks, *i.e.*, FM, AS, and SM tasks, CDND either matches or comes very close to the top-performing models, validating its status as a consistently high-performing model. In Appendix A.3, we have visualization results to highlight the qualitative differences.

Models	FA	FM	FS	MA	MF	MS	AF	AM	AS	SA	SF	SM	AVG
Supervised	$80.9_{\pm 7.2}$	$81.8_{\pm 0.3}$	$82.4_{\pm 1.2}$	$80.9_{\pm 7.2}$	$84.0_{\pm 1.8}$	$82.4_{\pm 1.2}$	$84.0_{\pm 1.8}$	$81.8_{\pm 0.3}$	$82.4_{\pm 1.2}$	$80.9_{\pm 7.2}$	$84.0_{\pm 1.8}$	$81.8_{\pm 0.3}$	82.3
Unsupervised	$78.5_{\pm 0.4}$	$60.9_{\pm 0.6}$	$66.5_{\pm 0.6}$	$26.6_{\pm 3.5}$	$33.6_{\pm 1.3}$	$69.9_{\pm 1.2}$	$38.5_{\pm 2.2}$	$31.2_{\pm 1.4}$	$30.0_{\pm 3.6}$	$74.1_{\pm 1.0}$	$68.4_{\pm 2.4}$	$64.5_{\pm 0.5}$	53.6
AdaptSegMap	$70.5_{\pm 3.4}$	$60.1_{\pm 0.6}$	$65.3_{\pm 1.3}$	$49.1_{\pm 9.7}$	$54.0_{\pm 0.5}$	$62.8_{\pm 7.6}$	$\mathbf{44.2}_{\pm 1.7}$	$35.4_{\pm 0.3}$	$35.1_{\pm 1.4}$	$70.1_{\pm 2.5}$	$67.7_{\pm 1.4}$	$63.8_{\pm 1.2}$	56.5
RS	$78.7_{\pm 0.5}$	$60.7_{\pm 0.4}$	$\mathbf{66.9}_{\pm 0.4}$	$59.6_{\pm 5.0}$	$38.4_{\pm 2.1}$	$\mathbf{70.4}_{\pm 1.0}$	$44.0_{\pm 0.6}$	$30.4_{\pm 0.5}$	$36.6_{\pm 0.8}$	$70.7_{\pm 0.8}$	$\mathbf{73.0}_{\pm 1.5}$	$\mathbf{65.3}_{\pm 1.3}$	57.9
DefRec+PCM	$78.8_{\pm 0.2}$	$\mathbf{60.9}_{\pm 0.8}$	$63.6_{\pm 0.1}$	$48.1_{\pm 0.4}$	$48.6_{\pm 2.4}$	$70.1_{\pm 0.8}$	$46.9_{\pm 1.0}$	$33.2_{\pm 0.3}$	$37.6_{\pm 0.1}$	$66.3_{\pm 1.7}$	$66.5_{\pm 1.0}$	$62.6_{\pm 0.2}$	56.9
GAST	$76.7_{\pm 2.3}$	$55.0_{\pm 1.0}$	$60.3_{\pm 1.0}$	$52.1_{\pm 4.4}$	$35.2_{\pm 0.4}$	$69.6_{\pm 1.2}$	$43.3_{\pm 3.7}$	$25.9_{\pm 3.6}$	$30.8_{\pm 4.0}$	$57.4_{\pm 10.6}$	$66.1_{\pm 1.3}$	$64.6_{\pm 0.5}$	53.1
ImplicitPCDA	$47.5_{\pm 0.6}$	$53.2_{\pm 1.0}$	$54.2_{\pm 3.4}$	$51.1_{\pm 1.6}$	$\mathbf{64.0}_{\pm 1.3}$	$56.1_{\pm 4.2}$	$44.1_{\pm 0.9}$	$\mathbf{42.3}_{\pm 1.3}$	$\mathbf{40.5}_{\pm 1.2}$	$49.7_{\pm 2.1}$	$70.6_{\pm 1.4}$	$55.0_{\pm 2.5}$	52.4
CDND	$\mathbf{81.5}_{\pm 2.0}$	$60.7_{\pm 0.5}$	$61.4_{\pm 0.5}$	$\mathbf{68.6}_{\pm 1.4}$	$47.2_{\pm 1.4}$	$67.7_{\pm 1.4}$	$43.6_{\pm 0.5}$	$35.3_{\pm 2.2}$	$40.1_{\pm 1.5}$	$\mathbf{77.5}_{\pm 0.5}$	$70.4_{\pm 1.1}$	$65.1_{\pm 0.3}$	**59.9**

Table 3: Performance results (mIOU) on PointSegDA dataset.

Overall, the performance improvement of our method, while consistent, is less pronounced compared to the results on PointDA-10. This is primarily due to the nature of the segmentation task, which involves fine-grained, per-point predictions. Our curvature diversity-driven deformation focuses on regional semantics and is more directly beneficial for object-level classification tasks. It is worth noting that other state-of-the-art methods also report smaller margins on PointSegDA. Nevertheless, our method still achieves the highest average mIoU, demonstrating its effectiveness in segmentation settings.

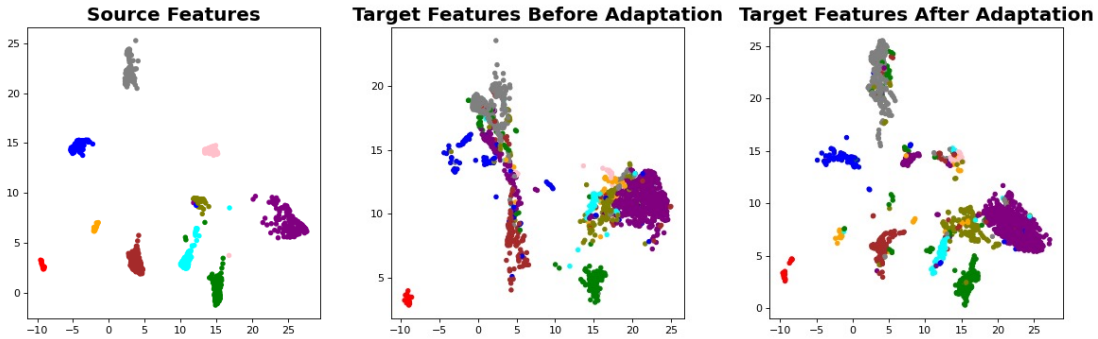

Figure 2: UMAP visualizations depict pre-activation data representations for the MS^+ task, with different colors denoting different classes. The center plot shows the target domain test data representations generated from a model trained on the source dataset without any adaptation. The left and right plots show the source and target domain data representations after adaptation using CDND.

5.4 Ablation Study

To demonstrate the effectiveness of each component of CDND, we conduct ablative studies on the PointDA-10 dataset. There are several ways to evaluate curvature diversity. While standard deviation is commonly used to evaluate the diversity of data points, we propose using entropy. We compare our entropy-based approach (CurvRec(En)) with a standard deviation-based method (CurvRec(S)). To validate our hypothesis that focusing on low curvature diversity regions can improve performance, we investigate the impact of deforming areas with both high (CurvRec(En)+High, CurvRec(S)+High) and low (CurvRec, CurvRec(S)+Low) curvature diversity.

Effectiveness of CurvRec. We compare against CurvRec(En) variants with CurvRec(S) variants in Table 2. We observe a distinct difference between CurvRec(En)-High and CurvRec. In contrast, there is a much less distinction between CurvRec(S)-High and CurvRec(S)-Low. This observation suggests that entropy is a superior method for evaluating curvature diversity in regions. Notably, all CurvRec measures outperform DefRec, regardless of whether the focus is on high or low curvature diversity. However, CurvRec outperforms CurvRec(En)-High, showing that is is more effective to deform low curvature regions. Compared to the plain NWD, all CurvRec variants perform better than plain NWD. Specifically, CurvRec surpasses

DefRec and NWD by approximately 3%. Though CurvRec demonstrates better performance overall, it does not outperform our proposed CDND. CDND (CurvRec +D-NWD), outperforms all CurvRec variants, DefRec variants, and plain NWD. This highlights the effectiveness of our D-NWD loss. Compared to CurvRec only, integrating with D-NWD improves average performance by 2.9%, with specific gains of 6.5% on MS^+ and 5.6% on SS^+.

Effectiveness of D-NWD. To illustrate the effectiveness of our D-NWD, we first compare CDND with two alternatives: CurvRec+PCM, which replaces D-NWD with PCM (PointMixup), and CurvRec+NWD. On average, CDND outperforms both methods. Specifically, compared to CurvRec+PCM, CDND achieves improvements of approximately 5% on MS^+, and 1% to 2% on SS^+, MS, and SM. When compared to CurvRec+NWD, CDND surpasses 4.4% on MS^+, 3% on SS^+, and 1.4% on S^+S. To further demonstrate the generalizability of D-NWD across different deformation methods, we include results for DefRec+D-NWD and DefRec+NWD. Compared to plain DefRec, DefRec+D-NWD achieves an overall improvement of 2.8%, with notable gains of 6.2% on MS^+, 3.4% on S^+S, and 8.3% on SS^+. Moreover, DefRec+D-NWD consistently outperforms DefRec+NWD on MS^+ and S^+S, as well as on average, although the margin is relatively small. Note that our theoretical analysis proves that D-NWD effectively reduces domain discrepancy across various deformation methods, but the degree to which it outperforms NWD depends on the deformation quality. With a well-designed deformations method like our proposed CurvRec, D-NWD can outperform NWD significantly. While D-NWD is crucial for tasks with large domain gaps (e.g., MS^+, S^+S, SS^+), its advantages are less significant in tasks with smaller domain gaps (e.g., MS, SM, S^+M), where CurvRec or DefRec alone already perform strongly. In these cases, plain CurvRec or DefRec performs better, so we retain their performance for MS, SM, and S^+M. Additionally, we test the use of $\mathcal{L}_{\text{NWD}}$ without adding $\mathcal{L}^{\mathcal{T}}_{\text{NWD}}$ to enhance consistency in the predictions of original target domain samples and deformed target domain samples, which corresponds to CDND with $\beta_2 = 0$. CDND with $\beta_2 = 0$ shows only a minor performance decrease compared to CDND, indicating that D-NWD is the primary contributor to good performance.

5.5 Analytic Experiments

We conduct analytical experiments to gain deeper insights into the effectiveness of our approach. Specifically, we assess how CDND impacts the *distribution* of the target domain in the classifier's output space for the challenging ModelNet to ScanNet task (MS^+); ModelNet is a synthetic dataset, while ScanNet is a real-world dataset, making the domain shift between them particularly challenging. We used UMAP to visualize and compare data representations of validation data from the source domain and test data from the target domain, both before and after applying CDND. Figure 2 shows each point as a data representation in the classifier's output space before softmax activation, with different colors denoting different classes. The middle plot in Figure 2 illustrates that, prior to adaptation, the classifier struggles with the target domain data, as points from different classes are heavily intermixed. However, after applying CDND, the class boundaries become more distinct, and the distribution of target domain representations aligns well with that of the source domain. This improvement is visible in the left and right plots of Figure 2, where the arrangement of points shows a more distinct pattern across both domains. In other words, we see that the feature space becomes domain-agnostic. This visualization demonstrates CDND's efficacy in reducing domain shift-induced performance degradation and enhancing class distinction. **More UMAP visualization analysis of other baselines can be found in Appendix A.3**.

6 Conclusion

We developed a new unsupervised domain adaptation approach for point cloud data. Our method integrates curvature diversity-based deformation with Deformation-based D-NWD to mitigate target domain performance degradation. Our theoretical analysis of D-NWD shows it minimizes an upper bound for target domain error. Additionally, we show that D-NWD can be applied to any deformation method. Experiments indicate that our approach surpasses SOTA methods on two point cloud benchmarks. Ablation studies confirm that both components of CDND are necessary for optimal performance. Future works include scenarios when source domain data is not accessible due to privacy or security concerns, or when the domains share only a subset of their classes.

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A Appendix

A.1 Proofs for Theorems

In this section, we first prove Theorem 1, which serves as the foundation for Theorem 2. Our proofs are structured as follows: we begin by proving Lemma 1, which supports a key assumption in Theorem 1. Next, we present the proof of Theorem 1. After proving Theorem 1, we prove Lemma 4 and conclude with the proof of Theorem 2.

Definition 2 (Probability Spacce). Durrett (2019) *A probability space is a triple $(\Omega, \mathcal{F}, \nu)$. Ω represents the sample space, the set of all possible outcomes. $\mathcal{F}$ represents the set of events and is a σ-algebra, which is a nonempty collection of subsets of Ω. $\mathcal{F}$ is closed under complements and countable unions. ν represents a probability measure on the measurable space $(\Omega, \mathcal{F})$. It is a function $\nu : \mathcal{F} \to [0, 1]$ that assigns to each event $A \in \mathcal{F}$ a real value $\nu(A)$ (the probability of A). ν satisfies the following three axioms:*

- *Non-negativity: For every $A \in \mathcal{F}$, $\nu(A) \geq v(\emptyset) = 0$.*
- *Normalization: $\nu(\Omega) = 1$.*
- *σ-additivity (Countable Additivity): For any countable sequence of pairwise disjoint events $A_1, A_2, A_3, \cdots \in \mathcal{F}$ (where $A_i \cap A_j = \emptyset$ for $i \neq j$),*

$$\nu\left(\bigcup_{i=1}^{\infty} A_i\right) = \sum_{i=1}^{\infty} \nu(A_i).$$

Lemma 1. *Let $(\Omega_1, \mathcal{F}_1, \nu_1)$ and $(\Omega_2, \mathcal{F}_2, \nu_2)$ be two probability spaces, where Ω_1, Ω_2 are two disjoint sample spaces. Let $p_1, p_2 \in [0, 1]$ be constants such that $p_1 + p_2 = 1$. Let $(\Omega_3, \mathcal{F}_3)$ be a measurable space, where $\mathcal{F}_3$ is the σ-algebra on $\Omega_3 = \Omega_1 \cup \Omega_2$. Then, the measure ν_3 defined on the measurable space $(\Omega_3, \mathcal{F}_3)$ as:*

$$\nu_3(A) = p_1\nu_1(A \cap \Omega_1) + p_2\nu_2(A \cap \Omega_2), \quad \forall A \in \mathcal{F}_3,$$

is a probability measure on $(\Omega_3, \mathcal{F}_3)$.

Proof. Since ν_1 and ν_2 are probability measures, they satisfy $\nu_1(B) \geq 0$ for all $B \in \mathcal{F}_1$ and $\nu_2(C) \geq 0$ for all $C \in \mathcal{F}_2$. For any set $A \in \mathcal{F}_3$, we have:

$$\nu_3(A) = p_1\nu_1(A \cap \Omega_1) + p_2\nu_2(A \cap \Omega_2).$$

Given that $p_1, p_2 \geq 0$ and $\nu_1(A \cap \Omega_1) \geq 0$ and $\nu_2(A \cap \Omega_2) \geq 0$, it follows that $\nu(A) \geq 0$. Thus, ν is non-negative. Then, we need to show that $\nu(\Omega_1 \cup \Omega_2) = 1$. Consider:

$$\nu_3(\Omega_1 \cup \Omega_2) = p_1\nu_1((\Omega_1 \cup \Omega_2) \cap \Omega_1) + p_2\nu_2((\Omega_1 \cup \Omega_2) \cap \Omega_2).$$

Since $(\Omega_1 \cup \Omega_2) \cap \Omega_1 = \Omega_1$ and $(\Omega_1 \cup \Omega_2) \cap \Omega_2 = \Omega_2$, and $\nu_1(\Omega_1) = 1$ and $\nu_2(\Omega_2) = 1$, we have:

$$\nu_3(\Omega_1 \cup \Omega_2) = p_1 \cdot 1 + p_2 \cdot 1 = p_1 + p_2 = 1.$$

Thus, ν_3 is normalized. Let $\{A_i\}_{i=1}^{\infty}$ be a countable collection of pairwise disjoint sets in $\mathcal{F}_3$. By definition of ν_3,

$$\nu_3\left(\bigcup_{i=1}^{\infty} A_i\right) = p_1\nu_1\left(\left(\bigcup_{i=1}^{\infty} A_i\right) \cap \Omega_1\right) + p_2\nu_2\left(\left(\bigcup_{i=1}^{\infty} A_i\right) \cap \Omega_2\right).$$

Since the A_i are pairwise disjoint, $(\bigcup_{i=1}^{\infty} A_i) \cap \Omega_1 = \bigcup_{i=1}^{\infty} (A_i \cap \Omega_1)$, and similarly for Ω_2. Using the σ-additivity of ν_1 and ν_2:

$$p_1\nu_1\left(\bigcup_{i=1}^{\infty}(A_i \cap \Omega_1)\right) = p_1 \sum_{i=1}^{\infty} \nu_1(A_i \cap \Omega_1),$$

$$p_2\nu_2\left(\bigcup_{i=1}^{\infty}(A_i\cap\Omega_2)\right)=p_2\sum_{i=1}^{\infty}\nu_2(A_i\cap\Omega_2).$$

Thus,

$$\nu_3\left(\bigcup_{i=1}^{\infty}A_i\right)=p_1\sum_{i=1}^{\infty}\nu_1(A_i\cap\Omega_1)+p_2\sum_{i=1}^{\infty}\nu_2(A_i\cap\Omega_2)=\sum_{i=1}^{\infty}\left(p_1\nu_1(A_i\cap\Omega_1)+p_2\nu_2(A_i\cap\Omega_2)\right).$$

Since $\nu_3(A_i)=p_1\nu_1(A_i\cap\Omega_1)+p_2\nu_2(A_i\cap\Omega_2)$, we get:

$$\nu_3\left(\bigcup_{i=1}^{\infty}A_i\right)=\sum_{i=1}^{\infty}\nu_3(A_i).$$

Thus, ν_3 satisfies σ-additivity. Since ν_3 satisfies non-negativity, normalization, and σ-additivity, by definition, ν_3 is a valid probability measure.
Now, for the following Lemmas and Theorems, we define:

$$\varepsilon_s(C_1,C_2)=\mathbb{E}_{f_s\sim\nu_s}\left[|C_1(f_s)-C_2(f_s)|\right],$$

$$\varepsilon_{s\cup s^d}(C_1,C_2)=\mathbb{E}_{\hat{f}_s\sim\nu_{s\cup s^d}}\left[\left|C_1(\hat{f}_s)-C_2(\hat{f}_s)\right|\right].$$

where C_1,C_2 are two classifiers. We define $\varepsilon_t(C_1,C_2)$ and $\varepsilon_{t\cup t^d}(C_1,C_2)$ in the same manner.

Lemma 2 (Lemma 1 Chen et al. (2022)). *Let ν,ν' be two probability measures on $(\Omega,\mathcal{F})$. Let $d(x,y)$ be the distance between $x\sim\nu$ and $y\sim\nu'$. W_N represents the NWD. Given a family of classifiers $C\in\mathcal{H}_1$ and a ideal classifier $C^*\in\mathcal{H}_1$ satisfying the K_L-Lipschitz constraint, where $\mathcal{H}_1$ is a subspace of $\mathcal{H}$, the following holds for every $C,C^*\in\mathcal{H}_1$.*

$$|\varepsilon(C,C^*)-\varepsilon'(C,C^*)|\leq 2K_L\cdot W_N(\nu_1,\nu_2),$$

where ε is the error on ν and ε' is the error on ν'.

Theorem 1. *Let $(\Omega_o,\mathcal{F}_o,\nu_s)$, $(\Omega_d,\mathcal{F}_d,\nu_{s^d})$, $(\Omega_o,\mathcal{F}_o,\nu_t)$, and $(\Omega_d,\mathcal{F}_d,\nu_{t^d})$ be four probability spaces, where Ω_o and Ω_d are disjoint and $\Omega_o\cup\Omega_d\subseteq\mathbb{R}^n$. With the results of Lemma 1, let $(\Omega_o\cup\Omega_d,\mathcal{F}_u,\nu_{s\cup s^d})$ and $(\Omega_o\cup\Omega_d,\mathcal{F}_u,\nu_{t\cup t^d})$ be two probability spaces with probability measures defined as $\nu_{s\cup s^d}=1/2\nu_s+1/2\nu_{s^d}$ and $\nu_{t\cup t^d}=1/2\nu_t+1/2\nu_{t^d}$. Specifically, when sampling from $\nu_{t\cup t^d}$, there is an equal probability of $1/2$ to sample from ν_t or ν_{t^d}. Similarly, sampling from $\nu_{s\cup s^d}$ gives an equal probability of $1/2$ to draw from ν_s or ν_{s^d}. Consider a classifier $C\in\mathcal{H}_1$ and an ideal classifier $C^*=\arg\min_C\varepsilon_{s\cup s^d}(C)+\varepsilon_t(C)$ satisfying the K_L-Lipschitz constraint, where $\mathcal{H}_1$ is a subspace of the hypothesis space $\mathcal{H}$. For every classifier C in $\mathcal{H}_1$, the following inequality holds:*

$$\varepsilon_t(C)\leq\ 2\varepsilon_{s\cup s^d}(C)+4K_L\cdot W_N(\nu_{s\cup s^d},\nu_{t\cup t^d})+\eta^*,$$

where $\eta^*=2\varepsilon_{s\cup s^d}(C^*)+\varepsilon_t(C^*)$ is the ideal combined error and is a sufficiently small constant.

Proof. Let Z be an indicator random variable that indicates whether the sample $\hat{f}_t$ is drawn from ν_t or ν_{t^d}:

- $Z=0$ if the sample is from ν_{t^d}.
- $Z=1$ if the sample is from ν_t.

By the Law of Total Expectation, we have:

$$\begin{aligned}\varepsilon_{t\cup t^d}(C,C^*)=&\ \mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)|]\\=&\ \mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)|\mid Z=0]P(Z=0)+\mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)|\mid Z=1]P(Z=1).\end{aligned}$$

Substituting $P(Z=0)=p_0$ and $P(Z=1)=p_1$,

$$\varepsilon_{t\cup t^d}(C,C^*) = p_0\mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)| \mid Z=0] + p_1\mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)| \mid Z=1].$$

Recognize that $\mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)| \mid Z=1]$ is the expectation when $\hat{f}_t$ is drawn from ν_t,

$$\varepsilon_t(C,C^*) = \mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)| \mid Z=1].$$

Combining these, we get:

$$\varepsilon_{t\cup t^d}(C,C^*) = p_0\mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)| \mid Z=0] + p_1\varepsilon_t(C,C^*).$$

Then,

$$\frac{1}{p_1}\varepsilon_{t\cup t^d}(C,C^*) = \frac{p_0}{p_1}\mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)| \mid Z=0] + \varepsilon_t(C,C^*).$$

Since $\frac{p_0}{p_1}\mathbb{E}_{\hat{f}_t\sim\nu_{t\cup t^d}}[|C(\hat{f}_t)-C^*(\hat{f}_t)| \mid Z=0] \geq 0$,

$$\frac{1}{p_1}\varepsilon_{t\cup t^d}(C,C^*) \geq \varepsilon_t(C,C^*).$$

Substituting $p_1 = 1/2$, we obtain:

$$2\varepsilon_{t\cup t^d}(C,C^*) \geq \varepsilon_t(C,C^*).$$

Based on Lemma 2, we have:

$$|\varepsilon_{s\cup s^d}(C,C^*) - \varepsilon_{t\cup t^d}(C,C^*)| \leq 2K_L \cdot W_N(\nu_{s\cup s^d}, \nu_{t\cup t^d}).$$

By triangular inequality,

$$\begin{aligned}\varepsilon_t(C) &\leq \varepsilon_t(C^*) + \varepsilon_t(C^*,C),\\ \varepsilon_{s\cup s^d}(C,C^*) &\leq \varepsilon_{s\cup s^d}(C) + \varepsilon_{s\cup s^d}(C^*).\end{aligned}$$

Then, we can derive:

$$\begin{aligned}\varepsilon_t(C) &\leq \varepsilon_t(C^*) + \varepsilon_t(C^*,C)\\ &\leq \varepsilon_t(C^*) + 2\varepsilon_{t\cup t^d}(C^*,C)\\ &= \varepsilon_t(C^*) + 2\varepsilon_{s\cup s^d}(C,C^*) + 2\varepsilon_{t\cup t^d}(C,C^*) - 2\varepsilon_{s\cup s^d}(C,C^*)\\ &\leq \varepsilon_t(C^*) + 2\varepsilon_{s\cup s^d}(C,C^*) + 4K_L \cdot W_N(\nu_{s\cup s^d}, \nu_{t\cup t^d})\\ &\leq \varepsilon_t(C^*) + 2\varepsilon_{s\cup s^d}(C) + 2\varepsilon_{s\cup s^d}(C^*) + 4K_L \cdot W_N(\nu_{s\cup s^d}, \nu_{t\cup t^d})\\ &= 2\varepsilon_{s\cup s^d}(C) + 4K_L \cdot W_N(\nu_{s\cup s^d}, \nu_{t\cup t^d}) + \eta^*.\end{aligned}$$

Definition 3 (L_1-Transportation Cost Information Inequality). Djellout et al. (2004) *Given $\eta > 0$, a probability measure ν on a measurable space $(\Omega, \mathcal{F})$ satisfies $T_1(\eta)$ if the inequality*

$$W_1(\nu', \nu) \leq \sqrt{\frac{2}{\eta}H(\nu'|\nu)}$$

where

$$H(\nu'|\nu) = \int \log\frac{d\nu'}{d\nu}d\nu'$$

holds for any probability measure ν' on $(\Omega, \mathcal{F})$, where W_1 represents the 1-Wasserstein distance.

Lemma 3. (Corollary 2.6 in Bolley & Villani (2005)) *For a probability measure ν on a measurable space $(\Omega, \mathcal{F})$, the following statements are equivalent:*

- *ν satisfies $T_1(\eta)$ inequality for some η that can be explicitly found.*
- *ν has a square-exponential moment, i.e., there exists $\alpha > 0$ such that*

$$\int_\Omega \exp(\alpha d(x,y)^2)\, d\nu(x) \text{ is finite}$$

for any $y \in \Omega$. Here, d is a measurable distance over Ω.

Lemma 4. *Let $(\Omega_1, \mathcal{F}_1, \nu_1)$ and $(\Omega_2, \mathcal{F}_2, \nu_2)$ be two probability spaces, where Ω_1 and Ω_2 are disjoint. Let $p_1, p_2 \in [0,1]$ be constants such that $p_1 + p_2 = 1$. Define a new measure ν_3 on a measurable space $(\Omega_3, \mathcal{F}_3)$, where $\Omega_3 = \Omega_1 \cup \Omega_2$:*

$$\nu_3(A) := p_1\nu_1(A \cap \Omega_1) + p_2\nu_2(A \cap \Omega_2), \quad \forall A \in \mathcal{F}_3.$$

Assume that ν_1 and ν_2 admit square-exponential moments: for some constants $\alpha_1, \alpha_2 > 0$, and for all $y_1 \in \Omega_1$, $y_2 \in \Omega_2$,

$$\int_{\Omega_1} \exp\left(\alpha_1 d_1(x,y_1)^2\right)\, d\nu_1(x) < \infty, \quad \int_{\Omega_2} \exp\left(\alpha_2 d_2(x,y_2)^2\right)\, d\nu_2(x) < \infty,$$

where $d_1 : \Omega_1 \times \Omega_1 \to [0,\infty)$ and $d_2 : \Omega_2 \times \Omega_2 \to [0,\infty)$ are distance functions. Then, ν_3 is a probability measure on $(\Omega_3, \mathcal{F}_3)$, and there exists a distance function $d : \Omega_3 \times \Omega_3 \to [0,\infty)$ and a constant $\alpha \in (0, \min\{\alpha_1, \alpha_2\})$ such that ν_3 admits a square-exponential moment: for all $y \in \Omega_3$,

$$\int_{\Omega_3} \exp\left(\alpha d(x,y)^2\right)\, d\nu_3(x) < \infty.$$

Proof. First, we define $d : \Omega_3 \times \Omega_3 \to [0,\infty)$:

$$d(x,y) = \begin{cases} d_1(x,y) & \text{if } x,y \in \Omega_1 \\ d_2(x,y) & \text{if } x,y \in \Omega_2 \\ C & \text{if } x \in \Omega_1, y \in \Omega_2 \text{ (or vice versa)} \end{cases}$$

where C is a finite constant chosen to ensure d is a metric on Ω_3. d can be expressed as:

$$d(x,y) = d_1(x,y)\mathbf{1}_{x,y\in\Omega_1} + d_2(x,y)\mathbf{1}_{x,y\in\Omega_2} + C\mathbf{1}_{x\in\Omega_1, y\in\Omega_2 \text{ or } x\in\Omega_2, y\in\Omega_1}$$

where $\mathbf{1}$ is the indicator function. Accodring to Theorem 1.9 (d) in Rudin (1987), the indicator functions are measurable because $\Omega_1 \times \Omega_1$, $\Omega_2 \times \Omega_2$, and $(\Omega_1 \times \Omega_2) \cup (\Omega_2 \times \Omega_1)$ are all measurable sets in $\Omega_3 \times \Omega_3$. d_1 and d_2 are also measurable functions by assumption. Therefore, d is a sum of products of measurable functions. Hence, d is measurable. For any y in Ω_1,

$$\begin{aligned}
&\int_{\Omega_3} \exp(\alpha d(x,y)^2)\, d\nu_3(x) \\
=& p_1 \int_{\Omega_1} \exp(\alpha d_1(x,y)^2)\, d\nu_1(x) + p_2 \int_{\Omega_2} \exp(\alpha d(x,y)^2)\, d\nu_2(x) \\
\leq& p_1 \int_{\Omega_1} \exp(\alpha_1 d_1(x,y)^2)\, d\nu_1(x) + p_2 \int_{\Omega_2} \exp(\alpha C^2)\, d\nu_2(x) \\
=& p_1 \int_{\Omega_1} \exp(\alpha_1 d_1(x,y)^2)\, d\nu_1(x) + p_2 \exp(\alpha C^2) < \infty
\end{aligned}$$

For any y in Ω_2,

$$\begin{aligned}
&\int_{\Omega_3} \exp(\alpha d(x,y)^2)\, d\nu_3(x) \\
=& p_1 \int_{\Omega_1} \exp(\alpha d(x,y)^2)\, d\nu_1(x) + p_2 \int_{\Omega_2} \exp(\alpha d_2(x,y)^2)\, d\nu_2(x) \\
\leq& p_1 \int_{\Omega_1} \exp(\alpha C^2)\, d\nu_1(x) + p_2 \int_{\Omega_2} \exp(\alpha_2 d_2(x,y)^2)\, d\nu_2(x) \\
=& p_1 \exp(\alpha C^2) + p_2 \int_{\Omega_2} \exp(\alpha_2 d_2(x,y)^2)\, d\nu_2(x) < \infty
\end{aligned}$$

This proves that ν_3 has a square-exponential moment for some $0 < \alpha < \min(\alpha_1, \alpha_2)$.

Lemma 5. (Theorem 1.1 of Bolley et al. (2007)) *Let ν be a probability measure on $(\Omega, \mathcal{F})$ where $\Omega \subseteq \mathbb{R}^n$. ν satisfies a $T_1(\eta)$ inequality. Let $\hat{\nu} = \frac{1}{N}\sum_{i=1}^N \delta_{f^i}$ be its associated empirical measure defined on a sample set $\{f^i\}_{i=1}^N$ of size N drawn i.i.d from ν. Then for any $n' > n$ and $\eta' < \eta$, there exists some constant N_0 depending on n' and some square-exponential moment of ν such that for any $\epsilon > 0$ and $N \geq N_0 \max(\epsilon^{-(n'+2)}, 1)$, the following holds:*

$$\mathbb{P}[W_N(\nu, \hat{\nu}) > \epsilon] \leq \exp\left(-\frac{\eta'}{2} N \epsilon^2\right).$$

Theorem 2. *Under the assumption of Theorem 1, Ω_o and Ω_d are disjoint and $\Omega_o \cup \Omega_d \subseteq \mathbb{R}^n$. Let $(\Omega_o \cup \Omega_d, \mathcal{F}_u, \nu_{s\cup s^d})$ and $(\Omega_o \cup \Omega_d, \mathcal{F}_u, \nu_{t\cup t^d})$ be two probability spaces with $\nu_{s\cup s^d} = 1/2\nu_s + 1/2\nu_{s^d}$ and $\nu_{t\cup t^d} = 1/2\nu_t + 1/2\nu_{t^d}$, where $\nu_s, \nu_{s^d}, \nu_t, \nu_{t^d}$ each has a square-exponential moment. From Lemma 3 and 4, $\nu_{s\cup s^d}$ satsifies $T_1(\eta_s)$ for some η_s and $\nu_{t\cup t^d}$ satsifies $T_1(\eta_t)$ for some η_t. Let $\{\hat{f}_s^i\}_{i=1}^{N_s}$ and $\{\hat{f}_t^i\}_{i=1}^{N_t}$ be two sample sets of size N_s and N_t drawn i.i.d from $\nu_{s\cup s^d}$ and $\nu_{t\cup t^d}$, respectively. $\hat{\nu}_{s\cup s^d} = \frac{1}{N_s}\sum_{i=1}^{N_s} \delta_{\hat{f}_s^i}$ and $\hat{\nu}_{t\cup t^d} = \frac{1}{N_t}\sum_{i=1}^{N_t} \delta_{\hat{f}_t^i}$ are associated empirical probability measures. Then, for any $n' > n$ and $\eta' < \min(\eta_s, \eta_t)$, there exists a constant N_0 depending on n' such that for any $\delta > 0$ and $\min(N_s, N_t) \geq N_0 \max(\delta^{-(n'+2)}, 1)$, with probability at least $1 - \delta$, the following holds for all C:*

$$\varepsilon_t(C) \leq\ 2\varepsilon_{s\cup s^d}(C) + 4K_L \cdot W_N(\hat{\nu}_{s\cup s^d}, \hat{\nu}_{t\cup t^d}) + \eta^* + 4K_L \cdot \sqrt{\frac{2}{\eta'} \log \frac{1}{\delta}} \left(\sqrt{\frac{1}{N_s}} + \sqrt{\frac{1}{N_t}}\right), \tag{17}$$

where $\eta^* = 2\varepsilon_{s\cup s^d}(C^*) + \varepsilon_t(C^*)$ is the ideal combined error and is a sufficiently small constant.

Proof. Based on Theorem 1,

$$\varepsilon_t(C) \leq\ 2\varepsilon_{s\cup s^d}(C) + 4K_L \cdot W_N(\nu_{s\cup s^d}, \nu_{t\cup t^d}) + \eta^*.$$

As a part of a broader class of Wasserstein distances, W_N satisfies the axioms of a distance Villani et al. (2009). Hence, W_N satisfies the triangle inequality:

$$\begin{aligned}
\varepsilon_t(C) \leq& 2\varepsilon_{s\cup s^d}(C) + 4K_L \cdot W_N(\nu_{s\cup s^d}, \hat{\nu}_{s\cup s^d}) + 4K_L \cdot W_N(\hat{\nu}_{s\cup s^d}, \nu_{t\cup t^d}) + \eta^* \\
\leq& 2\varepsilon_{s\cup s^d}(C) + 4K_L \cdot W_N(\nu_{s\cup s^d}, \hat{\nu}_{s\cup s^d}) + 4K_L \cdot W_N(\hat{\nu}_{s\cup s^d}, \hat{\nu}_{t\cup t^d}) + 4K_L \cdot W_N(\hat{\nu}_{t\cup t^d}, \nu_{t\cup t^d}) + \eta^*.
\end{aligned}$$

From Lemma 5,

$$W_N(\nu_{s\cup s^d}, \hat{\nu}_{s\cup s^d}) \leq \sqrt{\frac{2}{\eta'} \log\left(\frac{1}{\delta}\right)} \cdot \sqrt{\frac{1}{N_s}},$$

$$W_N(\nu_{t\cup t^d}, \hat{\nu}_{t\cup t^d}) \leq \sqrt{\frac{2}{\eta'} \log\left(\frac{1}{\delta}\right)} \cdot \sqrt{\frac{1}{N_t}}.$$

W_N belongs to the family of 1-Wasserstein distance. By the symmetry property of distance,

$$W_N(\hat{\nu}_{s\cup s^d}, \nu_{s\cup s^d}) = W_N(\nu_{s\cup s^d}, \hat{\nu}_{s\cup s^d}) \leq \sqrt{\frac{2}{\eta'} \log\left(\frac{1}{\delta}\right)} \cdot \sqrt{\frac{1}{N_s}},$$

$$W_N(\hat{\nu}_{t\cup t^d}, \nu_{t\cup t^d}) = W_N(\nu_{t\cup t^d}, \hat{\nu}_{t\cup t^d}) \leq \sqrt{\frac{2}{\eta'} \log\left(\frac{1}{\delta}\right)} \cdot \sqrt{\frac{1}{N_t}}.$$

Substituting back, we have:

$$\varepsilon_t(C) \leq 2\varepsilon_{s\cup s^d}(C) + 4K_L \cdot W_N(\hat{\nu}_{s\cup s^d}, \hat{\nu}_{t\cup t^d}) + \eta^* + 4K_L \cdot \sqrt{\frac{2}{\eta'} \log \frac{1}{\delta}} \left(\sqrt{\frac{1}{N_s}} + \sqrt{\frac{1}{N_t}}\right).$$

A.2 Implementation Details

Our code is based on the open-source implementation of the DefRec+PCM. We trained our three CDND models with seeds {1, 2, 3} on A100 GPUs. For the PointSegDA dataset, we fixed the learning rate to be 0.001 and conducted a grid search to optimize the hyperparameters α, γ, β_1, and β_2 for each task. The specific hyperparameter values can be found in Table 5. Similarly, for the PointDA dataset, the hyperparameters are listed in Table 4. Training on the PointDA dataset takes approximately 10 hours, and around 8 hours required for source-only training without any adaptation method. This results in a high computational cost for hyperparameter tuning. Therefore, we do not tune the hyperparameters extensively. Similarly, for GAST and ImplicitPCDA, we use the hyperparameters provided in their open-source code for the PointDA dataset.

However, GAST and ImplicitPCDA have not been tested on the PointSegDA dataset before. When implementing GAST, we conduct a grid search on the PointSegDA dataset, exploring values of 0.1, 0.2, 0.5, and 1.0 for both $\mathcal{L}_{\text{rot}}$ and $\mathcal{L}_{\text{loc}}$. For ImplicitPCDA, we perform a grid search on the PointSegDA dataset, considering values of 0.1, 0.2, 0.5, and 1.0 for $\mathcal{L}_M$. Please refer to the original papers for the definitions of $\mathcal{L}_{\text{rot}}$, $\mathcal{L}_{\text{loc}}$, and $\mathcal{L}_M$.

Hyperparameter	**Values**
Learning Rate	0.001, 0.0001 (S^+M, MS)
α	0.5
γ	0.5
β_1	[0.0, 1.0]
β_2	0.2
# of Epochs	200

Table 4: Hyperparameters for PointDA.

Hyperparameter	**Values**
Learning Rate	0.001
α	1.0
γ	[0.05, 0.1, 0.2, 0.5, 1.0]
β_1	[0.0, 0.05, 0.1, 0.2, 0.5, 1.0]
β_2	[0.0, 0.2]
# of Epochs	150

Table 5: Hyperparameters for PointSegDA.

For our method, we set the number of regions to be deformed to 5 for PointDA and 10 for PointSegDA. Each region contains 55 points. The total number of regions is set to 20 for the PointDA dataset and 40 for the PointSegDA dataset. For DefRec+PCM and CurvRec+PCM, we follow the hyperparameter settings used in the open-source implementation of DefRec and PCM (PointMixup).

Challenges of Applying SPST with mIoU The mIoU metric is defined as:

$$\text{mIoU} = \frac{1}{M} \sum_{m=1}^{M} \frac{TP_m}{TP_m + FP_m + FN_m}$$

$$\text{where:} \begin{cases} M = \text{number of classes} \\ TP_m = \text{true positive for class } m \\ FP_m = \text{false positive for class } m \\ FN_m = \text{false negative for class } m \end{cases}$$

SPST typically relies on ranking training samples by difficulty and gradually incorporating harder examples into training. The training samples for point cloud segmentation tasks are points in point clouds. However, mIoU is a global metric that evaluates performance across an entire point cloud, making it challenging to assign difficulty scores to individual points in a point cloud. The mechanism of SPST mismatches the per-point cloud, rather than per-point, evaluation criterion of mIoU.

A.3 Additional Results

We also added visualization comparisons on ModelNet to ScanNet task in Figure 3, featuring three methods: GAST and DefRec (which achieved second and third-best scores on the ModelNet to ScanNet task) and PCFEA (the most recently published method). Our approach achieves the most distinct and well-separated class clusters (represented by distinct colors) on the target domain test set, while PCFEA and GAST show considerable class mixing, and DefRec shows a tight clustering of some groups, lacking a clear separation.

We present segmentation visualization results in Figure 4 for the PointSegDA dataset. Different segmentation parts are highlighted using different colors. We selected three domains—MA, MF, and SA—to demonstrate the effectiveness of our CDND method, which achieves state-of-the-art performance with a significant margin over the second-best method. In the figure, the last row labeled "GT" shows the ground truth. Our CDND results are shown in the third row. Although CDND achieves 2% improvement in mIoU compared to RS Sauder & Sievers (2019), we include this comparison to highlight the qualitative differences in segmentation. The "Unsupervised" row represents results without any domain adaptation technique. From Figure 4, it is evident that CDND produces segmentation results most closely aligned with the ground truth, outperforming both RS and the unsupervised baseline. Additionally, It is worth noting that other methods also report small increases in mIoU margins on PointSegDA dataset.

We also conduct experiments to evaluate the effects of different Gaussian variances and the total number of regions in CurvRec. These experiments are performed on the FAUST to ADOBE task of the PointSegDA dataset, and the results are presented in Table 6. Specifically, we vary the total number of regions, $k \in \{10, 25, 40, 55, 70, 85, 100\}$, and the Gaussian variance, $\sigma^2 \in \{0.0001, 0.001, 0.01, 0.1\}$. The number of points per region, N_k, is calculated as $\lfloor \frac{2048}{k} \rfloor + 4$, where 2048 is the number of points per sample. The additional 4 points ensure that $k \times N_k \geq 2048$. This condition is important because we want to maximize coverage of the input point cloud—ideally assigning all 2048 points to regions. Without this constraint, $\lfloor \frac{2048}{k} \rfloor$ could be less than 2048, leaving some points uncovered. By slightly increasing the number of points per region, we ensure complete or near-complete coverage. As shown in the table, the optimal performance is achieved when the variance is set to 0.001 and the number of regions is 40, which aligns with our experimental settings. The results also indicate that model performance is relatively insensitive to the number of regions but is more sensitive to the choice of variance. In particular, large variances (e.g., 0.1) lead to more uniform deformations and may negatively affect performance.

$\sigma^2 \setminus (k, N_k)$	(10, 208)	(25, 85)	**(40, 55)**	(55, 41)	(70, 33)	(85, 28)	(100, 24)
0.0001	$80.5_{\pm 0.9}$	$80.3_{\pm 2.8}$	$79.8_{\pm 1.9}$	$79.6_{\pm 1.2}$	$79.4_{\pm 1.0}$	$80.6_{\pm 1.7}$	$80.1_{\pm 1.7}$
0.001	$80.8_{\pm 1.2}$	$80.7_{\pm 1.8}$	$\mathbf{81.5}_{\pm 2.0}$	$80.9_{\pm 1.7}$	$79.2_{\pm 2.3}$	$78.9_{\pm 2.1}$	$79.1_{\pm 2.0}$
0.01	$78.6_{\pm 2.9}$	$80.3_{\pm 2.3}$	$81.0_{\pm 1.9}$	$79.3_{\pm 4.0}$	$79.9_{\pm 2.3}$	$79.6_{\pm 3.0}$	$77.5_{\pm 4.0}$
0.1	$69.0_{\pm 2.8}$	$68.3_{\pm 3.6}$	$68.2_{\pm 5.6}$	$67.2_{\pm 4.4}$	$67.9_{\pm 4.9}$	$69.9_{\pm 4.5}$	$67.0_{\pm 5.2}$

Table 6: Results for varying σ^2, number of regions k, and points per region N_k.

Another important hyperparameter is the number of regions selected for deformation. In Table 7, we present results for the FAUST to ADOBE task of the PointSegDA dataset, where the total number of regions is fixed at $k = 40$ and the Gaussian variance is set to $\sigma^2 = 0.001$. We vary the number of deformed regions N (expressed as the ratio N/k) to observe its effect on performance. The results show that performance is not sensitive to the number of deformed regions. The optimal performance is achieved when the ratio is set to 0.25, which aligns with our experimental settings.

In the theoretical section, we note that the results can be generalized to arbitrary sampling ratios by adjusting the mixture weights in the underlying probability measures. While it is theoretically challenging to prove

Ratio (N/k)	0.10	**0.25**	0.30	0.45
FA	$80.3_{\pm 3.0}$	$\mathbf{81.5}_{\pm 2.0}$	$81.1_{\pm 2.3}$	$79.5_{\pm 3.3}$

Table 7: Results for varying the number of deformed regions N, expressed as the ratio N/k.

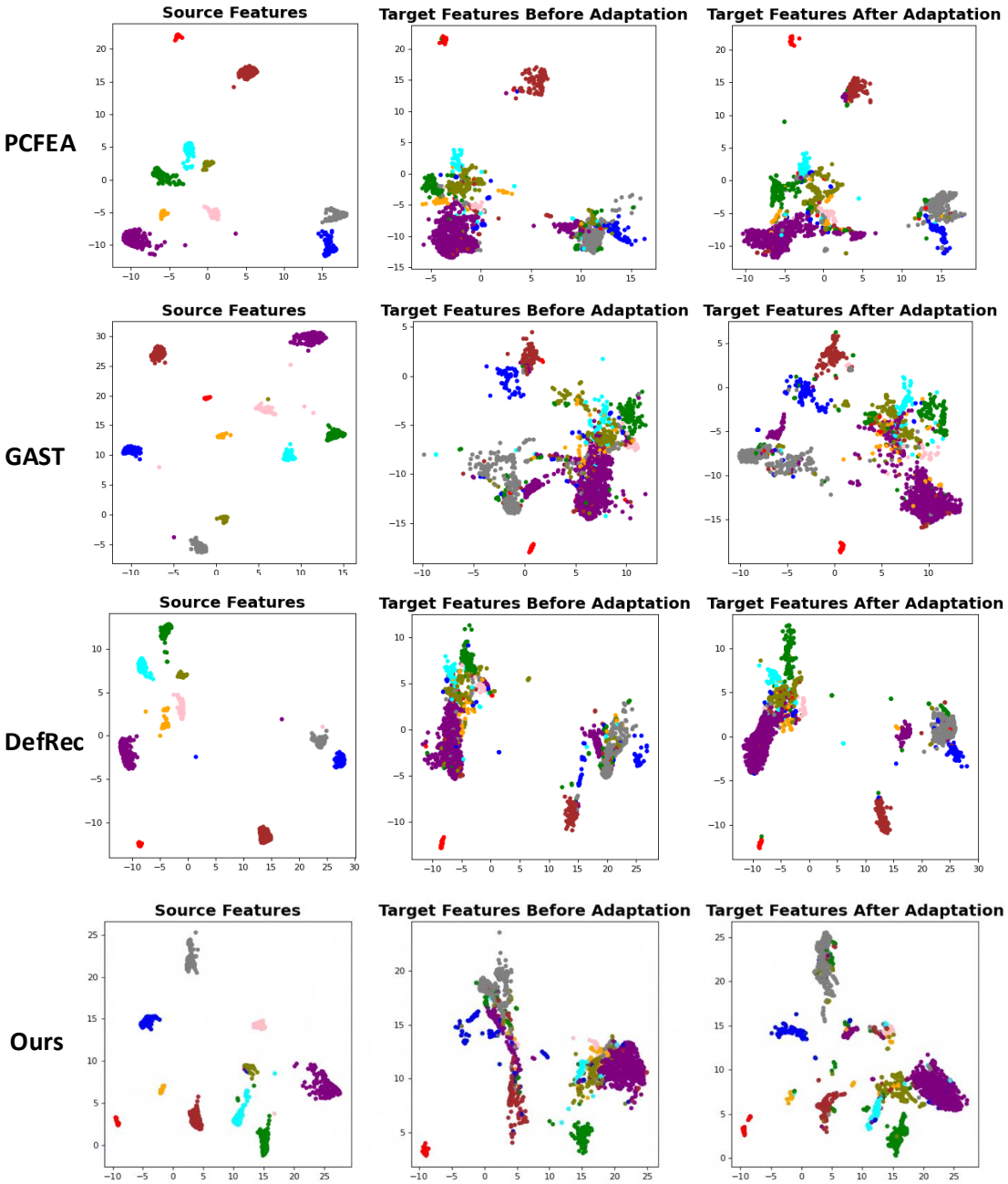

Figure 3: Umap Results of Other Baselines on PointDA-10 dataset.

that a specific sampling ratio between original and deformed samples (e.g., 1:1, 2:3, etc.) is universally optimal, our experimental results demonstrate that our chosen setting yields high performance. We conduct experiments on the FAUST to ABODE task of the PointSegDA dataset. As shown in Table 8, we achieve

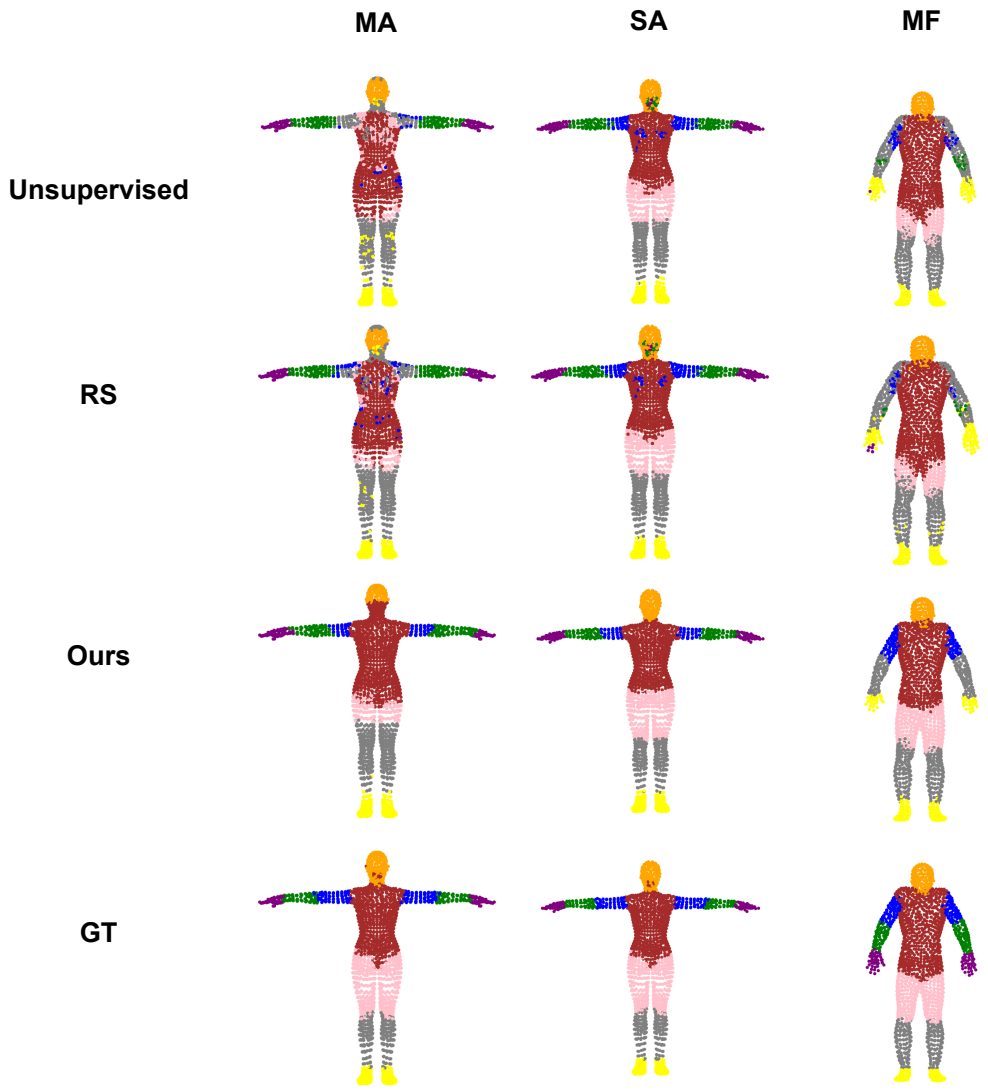

Figure 4: Segmentation Visualization Results of PointSegDA Dataset.

Ratio	1:1	1:2	1:3	1:4	2:1	2:3	2:4	3:1	3:2	3:4	4:1	4:2	4:3
FA	$81.5_{\pm 2.0}$	$80.1_{\pm 1.1}$	$79.3_{\pm 1.1}$	$79.5_{\pm 3.9}$	$79.5_{\pm 0.9}$	$79.3_{\pm 1.1}$	$79.5_{\pm 3.9}$	$76.7_{\pm 2.8}$	$79.9_{\pm 2.0}$	$79.5_{\pm 3.9}$	$77.2_{\pm 3.5}$	$77.1_{\pm 3.0}$	$\mathbf{81.7}_{\pm 1.2}$

Table 8: Results for various original:deformed sample ratios.

Neighborhood	10	20	40	CDND (32)
MS (Accuracy)	$\mathbf{84.1}_{\pm 0.3}$	$83.7_{\pm 0.4}$	$83.6_{\pm 0.4}$	$\mathbf{84.1}_{\pm 0.3}$
FA (mIOU)	$77.7_{\pm 2.9}$	$76.7_{\pm 3.5}$	$78.4_{\pm 2.6}$	$\mathbf{81.5}_{\pm 2.0}$

Table 9: Results for various neighborhoods.

Gaussian Variance	0.0001	0.01	0.1	Clean
MS (Accuracy)	$83.2_{\pm 0.5}$	$83.5_{\pm 0.4}$	$82.7_{\pm 1.5}$	$\mathbf{83.6}_{\pm 0.4}$
FA (mIOU)	$79.9_{\pm 1.3}$	$75.8_{\pm 5.1}$	$56.4_{\pm 1.9}$	$\mathbf{78.4}_{\pm 2.6}$

Table 10: Results for adding Gaussian noise.

the highest mIoU with the 1:1 and 4:3 ratios. This supports the intuition that a balanced and sufficiently diverse mix of original and deformed samples helps the model generalize more effectively across domains. It is also worth noting that our method is not highly sensitive to changes in the sampling ratio.

We also conducted additional experiments to analyze the impact of hyperparameters on curvature estimation. Curvature estimation depends on clean local geometry and can be influenced by real-world scanning artifacts. To evaluate, we report classification accuracy on the ModelNet to ShapeNet task from the PointDA dataset, and segmentation performance (mIOU) on the FAUST to ADOBE task from the PointSegDA dataset. Table 9 presents the results of varying the number of neighbors used in curvature computation. The performance remains relatively stable (though the segmentation task is more sensitive) across different neighborhood sizes, indicating that curvature estimation is not significantly sensitive to this parameter. Notably, our experiment setting of 32 neighbors (labeled as CDND) yields the best overall performance for both tasks. Regarding the number of principal components, we do not arbitrarily choose how many to use — the number is determined by the dimensionality of the input point cloud data. In our case, the input points lie in 3D space, so the covariance matrix is 3×3, yielding three eigenvalues and corresponding eigenvectors.

In Table 10, we introduce Gaussian noise to the point coordinates to evaluate the robustness of curvature estimation. For these experiments, the number of neighbors is fixed at 40. The classification task (ModelNet to ShapeNet) demonstrates robustness to noise, as the accuracy remains relatively consistent even with increasing noise variance. However, the segmentation task (FAUST to ADOBE) is more sensitive—its performance degrades significantly with higher noise levels. This increased sensitivity in segmentation arises because segmentation tasks require more precise, point-level geometric details to assign accurate labels to each point. Noise distorts these fine-grained features, making it harder for the model to distinguish between regions, especially in complex structures. In contrast, classification tasks benefit from more global shape features and can tolerate localized noise to a greater extent.

Method	Unsupervised	DANN	PointDAN	RS	DefRec + PCM	PCFEA	GAST	ImplicitPCDA	CDND
Time (hours)	8.0	9.0	9.2	35.9	10.4	8.3	18.9	9.8	10.4

Table 11: Training time (in hours) of different methods.

Table 11 presents the training time (in hours) for various baseline methods along with our proposed approach, CDND. As observed, our method does not introduce a significant computational overhead compared to the "Unsupervised" baseline (without any domain adaptation method), requiring extra 2.4 additional hours. This demonstrates a decent efficiency of CDND, especially when contrasted with methods like RS and GAST, which have significantly higher training costs. Furthermore, CDND's training time is still comparable to DANN, PointDAN, and DefRec+PCM. Overall, CDND is practically viable without significantly sacrificing efficiency.