On the Robustness of Average Losses for Partial-Label Learning

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Abstract—Partial-label learning (PLL) utilizes instances with PLs, where a PL includes several candidate labels but only one is the true label (TL). In PLL, identification-based strategy (IBS) purifies each PL on the fly to select the (most likely) TL for training; average-based strategy (ABS) treats all candidate labels equally for training and let trained models be able to predict TL. Although PLL research has focused on IBS for better performance, ABS is also worthy of study since modern IBS behaves like ABS in the beginning of training to prepare for PL purification and TL selection. In this paper, we analyze why ABS was unsatisfactory and propose how to improve it. Theoretically, we propose two problem settings of PLL and prove that average PL losses (APLLs) with bounded multi-class losses are always robust, while APLLs with unbounded losses may be non-robust, which is the first robustness analysis for PLL. Experimentally, we have two promising findings: ABS using bounded losses can match/exceed state-of-the-art performance of IBS using unbounded losses; after using robust APLLs to warm start, IBS can further improve upon itself. Our work draws attention to ABS research, which can in turn boost IBS and push forward the whole PLL.

Index Terms—Partial-label learning, robust loss, robustness analysis, weakly supervised learning.

I. INTRODUCTION

Deep neural networks (DNNs) have become the par excellence model in diverse application domains, which transform the input data (e.g., images) to the specific outputs (e.g., classes). Much of the success in running DNNs is attributed to its internal capability to approximate arbitrarily complex functions mapping input to output [1], [2], [3], [4], as well as an external driving force—labeled training data. It is widely believed that the performance of DNNs is improved as the number of data increases, reaching saturation only when millions of data are available [43], [54], [57], [77]. Their remarkable performance usually comes at a prohibitively high labeling cost, especially when data labeling must be carried out professionally. A shortage of skilled experts, an expensive and time-consuming labeling process, and privacy issues can pose challenges to the acquisition of high-quality labels. As a result, learning with imperfect but inexpensive labels is practically significant.

Crowdsourcing [7] relying on non-expert workers has recently emerged as an attractive surrogate. Unlabeled instances are typically assigned to workers of varying knowledge, and limited by their expertise, they often have difficulty recognizing the exact label from multiple ambiguous categories. Therefore, crowdsourcing platforms naturally allow workers to select several possible labels if they are uncertain about an instance. In this way, an instance is associated with a set of candidate labels where a fixed but unknown candidate is the true label. A set of candidate labels is referred to as a partial label (PL) for an instance, and the learning paradigm that can handle PLs is termed as partial-label learning (PLL) [8], [12], [18], [41], [42], [51], [61], [62], [66], [68], [74], [75], also known as ambiguous label learning [9], [10], [22], [67] and superset learning [27], [36], [37]. PLL attempts to infer the optimal multi-class classifier that is able to accurately predict the true label for unseen instances by fitting
PLs, and more ideally, the hypotheses can be modeled by DNNs. PLL problems arise in real-world scenarios [36], [40], [70] as well.

Research on PLL dates back about 20 years. Initially, Jin and Ghahramani [30] built up a maximum likelihood model to reassign class-posterior probabilities to candidate labels iteratively. This work opened up a main research route of PLL that purifies each PL on the fly to select the most likely true label during training [10], [16], [17], [30], [61], [66], which is named the identification-based strategy (IBS). Because IBS aims at eliminating the ambiguity [74] between individual instances and their true labels in the training phase, this technique is also commonly known as disambiguating [12], [30]. Contrariwise, Hüllermeier and Beringer [26] formalized PLL as a collaborative problem, where all candidate labels contribute to the learning objective equally. The idea is that the inductive bias underlying the learning process can benefit disambiguating the given PLs, and let trained models be able to predict the true label of any instance. Such a scheme is called the average-based strategy (ABS) [12], [73].

In recent years, the research of PLL has focused on IBS, since it was believed that the performance of IBS is more promising, while little attention has been paid to ABS. This is especially so in the era of deep learning [18], [41], [61]. The pessimism about ABS comes from “memorization” of over-parameterized DNNs, that is, the perfectly fitted DNNs can memorize all training samples, even if their labels are completely arbitrary [15], [71]. ABS is free from identifying the latent true labels during training, and therefore memorize all candidate labels. Then the PLL problem would be degenerated to a multiple-label problem [30], where it is acceptable that an arbitrary candidate label is taken for the “pseudo true label”. Then will ABS fail in the true-label prediction and should ABS just be phased out by the times?

In this paper, we argue the research value of ABS by showing its practical potential and theoretical superiority, problematizing the traditional view of ABS and pushing forward PLL as a whole. 

**Promising experimental findings.** Our work is inspired by a set of exploratory experiments. We propose a family of ABS losses named average partial-label (APL) losses, which are defined as the average of multi-class losses over all candidate labels. The categorical cross entropy (CCE) loss is the most popular multi-class loss in deep learning nowadays, and we observed that all existing deep IBS methods [18], [41], [61], [67] also adopted the CCE loss. Thus first, we trained several standard deep models with the APL loss equipped with the CCE loss on benchmark datasets where true labels were manually corrupted to PLs. In this case, we found that our ABS method with the CCE loss performs poorly as was previously thought, shown in Fig. 1. Extending these experiments, we then replaced the CCE loss with another widely-used loss, i.e., the generalized cross entropy (GCE) loss [76], and conducted experiments with early stopping [49]. Surprisingly, we observed that our ABS method with the GCE loss can often be on a par with, or even outperform a SOTA IBS method PRODEN [41], regardless of datasets, models, and optimizers. For the details of data generation processes, please see Section VI-A. These observations challenge common beliefs since one would expect that ABS is incapable of distinguishing true labels, thereby hurting generalization, while the results showed that ABS method with the APL loss can also predict the true label. Therefore, we would like to analyze what is it that distinguishes ABS methods that perform well from those that do not, and the answer to this question will hopefully help improve ABS.
To answer this question, we analyze the robustness of our ABS method to PLs, namely, whether the classification error on supervised data of the minimizer of the risk w.r.t. the APL losses is approximated to that of the Bayes classifier (learned using supervised data) [20], [21], [44], [46]. Thanks to the concise form of the APL losses, it is easy to estimate the risk under the APL losses from PLs and carry out empirical risk minimization (ERM). Thus we can analyze the robustness through existing mathematical techniques.

Furthermore, we take one step forward—unreliable PLL, which learns from noisy PLs, that is, the candidate-label set that might not include the true label. As the acquisition of training data expands, noise is inevitable, therefore, it is not an overstatement to say unreliable PLL is imminent in real-world applications. Unfortunately, previous PLL algorithms concentrate on noise-free PLs, and they have not been able to handle the unreliable PLL well, as shown in the rightmost columns in Fig. 1. To avoid confusion, we refer to the traditional PLL paradigm reliable PLL, to reliable PLL and unreliable PLL collectively as “PLL”, and to noise-free PL and noisy PL collectively as “PLL” in later sections.

To theoretically analyze the cause of success or failure of an ABS method, we formalize two problem settings for the generation processes of PLs, each with several specific instances that differ significantly in their conceptual approaches. With the help of them, we delve into multiple widely-used multi-class loss functions, and formally prove that APL losses with bounded loss functions (e.g., GCE) are always robust under mild assumptions on the domination of true labels, while APL losses with unbounded loss functions (e.g., CCE) may not be robust. The theoretical results are reconciled with experimental observations in Fig. 1. Given that there exists no such analysis for IBS yet, our robustness analysis is novel for not only ABS but also PLL.

ABS improvement to IBS. Moreover, we rethink the existing deep IBS methods. We point out that all modern IBS methods behave like ABS in the beginning of training to prepare for PL purification and true-label selection. In other words, they need to use ABS to warm start model training, and use the pretrained model to identify the true label. Thus they will select the correct true labels if ABS can become better, while they have hitherto used unbounded losses throughout the training process. As a consequence, IBS methods can in turn improve upon themselves by our study on ABS: we suggest utilizing bounded losses as a warmup for the first few epochs. We conduct extensive experiments to verify the effectiveness of this improvement.

Contributions. Our contributions can be summarized:

- We establish a theoretically grounded framework for ABS based on a simple yet effective APL loss family, the risk minimization of which is guaranteed to be robust under two problem settings for the data generation processes.
- To the best of our knowledge, we are the first to propose the unreliable PLL paradigm, further developing the practical potentiality of PLL in society. Our ABS method with the APL losses provides an effective baseline for unreliable PLL, and it also works well for reliable PLL without any modifications.
- We redraw the attention of the PLL community to ABS. Our research findings can not only improve ABS, but also enlighten a general principle to incorporate ABS into IBS methods to further enhance their performance, and push forward PLL as a whole.

Paper organization. We recapitulate the related work and discuss the philosophies behind ABS and IBS in Section II. In Section III, we give an overview of the problem settings and introduce the importance of robustness analysis. In Section IV, we propose APL losses and formalize generation processes of PLs. We present our main theoretical results and experimental findings in Section V and Section VI, respectively. We conclude in Section VII, and defer additional experimental results and all the proofs to the appendix, available online.

II. RELATED WORK

In this section, we review some seminal work in reliable PLL, discuss philosophies behind different technical routes of PLL, and analyze its relation to and difference from other machine learning problems.

Practical reliable PLL. [30] is one of the milestones for reliable PLL. It proposed to disambiguate noise-free PLs by using the expectation-maximization (EM) algorithm. In the E-step, the class-posterior probability is estimated as the normalization of current model predictions. In the M-step, the model parameters are updated in order to minimize the KL divergence between the given estimated probabilities and the model-based distributions. Such a strategy that identifies true labels along with model training is referred to as IBS. Following the milestone work, many IBS algorithms have been developed (e.g., [10], [17], [22], [42], [58], [66], [73], [75]).

Reliable PLL has also been studied along the other research route called ABS, pioneered by Hüllemeyer and Beringer [26]. They determined the class label for an unseen instance by voting among the candidate labels of its nearest neighbors. Cour et al. [12] proposed a convex loss that distinguishes the averaged output over the candidate labels from outputs over non-candidate labels. It is always believed that ABS is likely to fail as the outputs of pseudo true labels would overwhelm the output of true label. Therefore, the development of the two strategies was not balanced – IBS has been the focus of considerable recent research whilst ABS faded away.

From 2020, deep learning starts injecting new vitality into IBS. Almost at the same time, three works proposed to model classifiers by DNNs. Yao et al. [67] adopted ResNet as the backbone together with two specially designed regularizers for partially-labeled image classification. Yao et al. [68] used the co-teaching scheme [24] to let two networks interact with each other regarding the confidence levels of the instances. The method proposed by Lv et al. [41] progressively identifies true labels based on the memorization effect of DNNs [3], which is flexible on the learning models and loss functions. Later, Feng et al. [18] formalized for the first time the generation process of noise-free PLs, based on which they derived two provably consistent algorithms. Wen et al. [61] proposed a leveraged weighted loss to trade off the losses on candidate labels and...
non-candidate ones. Wu and Sugiyama [63] proposed a unified framework includes [18] as a special case. These three works are both compatible with DNNs.

**Provable reliable PLL.** Although the above practical algorithms have proven empirically successful on specific domains, there is an elusive theoretical gap in the understanding of them. Through the lens of learning theory, some researchers proposed seminal theoretical works in reliable PLL. Liu and Dietterich [37] proposed the small ambiguity degree condition to ensure that classification errors on any instance have a probability of being detected. The proof of this theorem requires strict assumptions: the approximation error equals zero and meanwhile the Bayes error equals zero (i.e., the deterministic scenario [48]). Cour et al. [12], Feng et al. [18], and Wen et al. [61] focused on the statistical consistency. They proposed a consistent loss based on some specific data generation process or deterministic scenario assumption, while our findings are general enough to hold under different generation processes and also a stochastic scenario [48].

**Philosophies behind IBS and ABS.** IBS iterates between the optimization of a learning model and the identification of the true label. Typically the identified true label has the biggest posterior of all labels, and must be in the candidate-label set. In other words, the "true" one is the "ideal" one. It implies that the true label can be uniquely determined given an input—it is satisfied only in the deterministic scenario where the class-posterior probability of the true label is equal to 1.

However, the natural world is more like the stochastic scenario that possesses some inherent randomness. In this setting, the label is a probabilistic function of the input, indicating that the same input will lead to an ensemble of unfixed output labels. ABS essentially gets rid of the deterministic scenario by avoiding recognizing the "ideal" label. The "true" label is considered as an "actually sampled" outcome, and consequently, the philosophy of ABS is compatible with the stochastic scenario. Therefore, it is crucial to design advanced methods and provide theoretical understandings for ABS.

**Relevant learning problems.** There are some weakly supervised learning problems related to PLL.

*Complementary-label (CL) learning* [19], [28], [29], [69] learns from weakly-supervised datasets wherein an instance is equipped with a CL. A CL specifies a class that the pattern does NOT belong to, so it can be considered as an extreme noise-free PL case with a fixed number \((k - 1)\) of candidate labels. Then from the algorithmic point, reliable PLL algorithms can directly handle the CL learning problem, but not also the other way around.

*Semi-supervised learning (SSL)* [6], [47], [59], [72], [78] learns from datasets consisting of both labeled and unlabeled data. Since we can regard the universe set of labels as the candidate labels of unlabeled data, SSL has some relation with PLL. However, standard SSL assumes that labeled data are fully supervised, which is different from reliable PLL, where labeled data are still ambiguous.

*Noisy-label learning (NLL)* [23], [24], [38], [50], [53] learns from noisy supervision where the training data are sampled from a corrupted distribution. Both NLL and PLL should have an underlying transition matrix linking the clean class posterior and the observed class posterior of an instance. Nonetheless, their matrix dimensions are different: the transition matrix is \(k \times k\) in NLL and \(k \times (2^k - 2)\) in PLL.

**III. PRELIMINARIES**

In this section, we formally introduce reliable PLL and propose unreliable PLL, and give the definition of robustness.

**A. Problem Setup**

**Basic Settings.** Let us consider a multi-class classification problem of \(k\) classes. Let \(\mathcal{X} \subseteq \mathbb{R}^d\) be the feature space, \(\mathcal{Y} = [k] \equiv \{1, 2, \ldots, k\}\) be the label space, and \(\mathcal{S} \equiv \{2^k \setminus \emptyset, \emptyset\}\) be the PL space. \(2^k\) means the collection of all subsets in \([k]\), and \(|\mathcal{S}| = 2^k - 2\) because the empty set and the whole label set are excluded. We denote by \(p(x, y)\) some probability density of “clean” distribution over \(\mathcal{X} \times \mathcal{Y}\). In fully-supervised classification, the goal is a learning model (e.g., a DNN) \(f : \mathbb{R}^d \rightarrow [k]\) that can make correct prediction on unseen inputs, with a set of i.i.d. supervised training data \(\{(x_i, y_i)\}_{i=1}^n\) sampled from \(p(x, y)\). A classifier \(f(x)\) is routinely assumed to take the following form:

\[
\hat{f}(x) = \arg \max_{i \in \mathcal{Y}} g_i(x),
\]

where \(g_i(x) : \mathbb{R}^d \rightarrow \mathbb{R}\) outputs a score for class \(i\). In this paper, we concentrate on deep learning: assume the learning model \(f\) is a DNN and apply softmax to convert scores into a vector of class-posterior probabilities, i.e., \(g_i(x) = p(i|\mathcal{X}) \in \Delta^{k-1}\) [55], where \(\Delta^{k-1}\) denotes the \(k\)-dimensional simplex.

While in PLL, for the notional clean distribution with probability density \(p(x, y)\), we instead observe i.i.d. PL training data \(\{(x_i, y_i, s_i)\}_{i=1}^n\) from a corrupted version \(p(x, s)\) of \(p(x, y)\) over \(\mathcal{X} \times \mathcal{S}\). The distribution \(p(x, s)\) is such that the marginal distribution of instances \(p(x)\) is unchanged, but the observed label is corrupted to an ambiguous candidate-label set. PLL tries to nonetheless learn the optimal classifier by fitting \(\{(x_i, s_i)\}_{i=1}^n\).

The key assumption in reliable PLL is that the PLs are noise-free, which means the latent true label \(y_i\) of an instance \(x_i\) is always included in its candidate-label set \(s_i\), i.e.,

\[
p(y_i \in s_i | x_i, s_i) = 1, \forall (x_i, y_i) \in p(x, y), \forall s_i \in \mathcal{S}.
\]

We argue that this assumption is fairly strict since the density \(p(x, y)\) of clean distribution is agnostic. For example, annotations of large-scale datasets usually need to be done by distributed workers in crowdsourcing platforms. Requiring crowdsourcing workers to cautiously judge each category to ensure that the correct one must be chosen partially runs counter to the original purpose of reducing labeling costs, and workers are unable to annotate tasks accurately due to the limited knowledge, even when combining labels provided by multiple workers, 100% accuracy is not guaranteed. Another vivid example is medical diagnosis. It is nontrivial for doctors to make an exact judgment on challenging cases with one single exam. Instead, giving several suspected diseases and referring patients to receiving other tests based on those suspicions are less demanding. However, owing to high inter- and intra-observer variability [23],
some complicated diseases are less noticeable and often difficult
to detect, even for experts, leading to unreliable candidate labels.
As the acquisition of training data expands, it is pervasive for
label information to be corrupted, but unfortunately, it has never
been considered in previous PLL works. Thus we introduce a
more general data setting titled unreliable PLL:

Definition 1 (unreliable PLL). Given the joint density
$p(x, y, s)$ and its marginal density $p(x, s)$, for any noisy PL
data $(x_i, s_i)$ independently sampled from $p(x, s)$, its true label
$y_i$ has a probability of $0 \leq \gamma \leq 1$ not being included in the
candidate-label set $s_i$, i.e.,

$$p(y_i \in s_i | x_i, s_i) = 1 - \gamma, \forall (x_i, y_i) \sim p(x, y), \forall s_i \in S,$$
where $\gamma$ is called the unreliability rate. Learning from noisy PL
data is called unreliable PLL.

B. Robustness

The $\ell$-risk of $f$ in fully-supervised learning w.r.t. multi-class
loss $\ell : \mathbb{R}^k \times \mathcal{Y} \to \mathbb{R}^+$ is defined as follows:

$$\mathcal{R}(f; \ell) = \mathbb{E}_{p(x,y)}[\ell(f(x), y)].$$

$\mathbb{E}$ denotes the expectation and its subscript indicates the dis-
tribution with respect to which the expectation is taken. The
Bayes optimal classifier that minimizes $\mathcal{R}(f; \ell_0)$ is given by
$f^* = \arg \min_f \mathcal{R}(f; \ell_0)$, where the optimality is defined over
all measurable functions. We denote by $\mathcal{R}^+ = \mathcal{R}(f^*)$ the cor-
responding Bayes risk under the clean distribution. Typically, $\ell$
is classification-calibrated [4], that is, the global minimizer of
$\mathcal{R}(f; \ell)$ is the same as that of $\mathcal{R}(f; \ell_0)$, which can be interpreted
as controlling the excess 0-1 risk by controlling the excess
$\ell$-risk [50].

Denote by $\tilde{\ell} : \mathbb{R}^k \times S \to \mathbb{R}^+$ a suitably modified $\ell$ for use
with PLs (defined in Section IV-A). Similarly, the PLL risk under
$p(x, s)$ w.r.t. PLL loss $\ell$ is defined as

$$\tilde{\mathcal{R}}(f; \tilde{\ell}) = \mathbb{E}_{p(x,s)}[\tilde{\ell}(f(x), s)].$$
The aim of PLL is to predict the true label for unseen instances.
However, most of the standard learning methods are hard to
perform well as they tend to exhibit overfitting on the candidate
labels in such scenarios [41].

Constructing robust losses from the perspective of the objec-
tive function is a powerful means in weakly supervised learning
[21, 44, 46]. Its focus is to derive the theoretical guarantee for
robust losses so that the learned classifier based on weak su-
pervision approximates the Bayes optimal classifier. Concretely,
a loss $\ell$ is robust to PLs (more specifically the risk minimization
with $\ell$ is asymptotically robust to PLs) if it guarantees that the
optimal PLL classifier $\hat{f}^* = \arg \min_f \tilde{\mathcal{R}}(f; \tilde{\ell})$ converges to the
Bayes optimal classifier.

Definition 2 (PL-robustness). We say that a loss $\tilde{\ell}$ is robust
and data (PL-robust) if for any $p(x, y, s) \in \mathcal{X} \times \mathcal{Y}$, $\mathcal{R}(f^*) - \mathcal{R}^*$
is bounded.

$$\mathcal{R}(f^*) - \mathcal{R}^*$$
is bounded means that $\hat{f}^*$ learned from PL data has a similar classification error to $f^*$ on the supervised data, i.e.,
minimizing $\tilde{\mathcal{R}}$ yields an approximate solution that minimizes $\mathcal{R}$. A guarantee of robustness thus sets an analogous calibration
theory [4] of PLL. Let $\hat{\mathcal{R}}^* = \mathcal{R}(\hat{f}^*)$. Then the robustness
condition will often be rewritten as that $\hat{\mathcal{R}}(f^*) - \hat{\mathcal{R}}^*$ is bounded,
which is slightly weak because it only signifies that $f^*$ is the
approximated minimizer of $\hat{\mathcal{R}}(f^*)$, but does not guarantee the
classification performance of $\hat{f}^*$ on the supervised data.

In statistical learning theory, consistency [48] is another
important concept. We use the superscript $*$ to indicate the
optimal solution over a given hypothesis class $\mathcal{F}$, i.e., $f^* = \arg \min_{f \in \mathcal{F}} \mathcal{R}(f; \ell)$. Suppose $\hat{f} = \arg \min_{f \in \mathcal{F}} \frac{1}{n} \sum_{i=1}^n \tilde{\ell}(f(x_i), s_i)$ is the PLL empirical risk mini-
mizer. The quality of $\hat{f}$ with respect to $f^*$ is measured by the
estimation error:

$$\mathcal{R}(\hat{f}) - \mathcal{R}(f^*) = (\mathcal{R}(\hat{f}) - \mathcal{R}(f^*) + (\mathcal{R}(f^*) - \mathcal{R}(f^*)).$$

If as $n \to \infty$, there is $\mathcal{R}(\hat{f}) \to \mathcal{R}(f^*)$, we say the PLL is con-
sistent. According to the universal approximation theorem [2],
[13] that using a proper DNN, the hypothesis space $\mathcal{F}$ is suffi-
ciently complex to contain the Bayes optimal classifier, we have
$f^* = f^*, f^* = f^*$. Thanks to this, the concepts of robustness
and consistency can be well connected in deep learning: RHS2 in
(1) is just the robustness measure. Therefore, consistency is a
sufficient but not a necessary condition of robustness. Although
robustness is a weaker property than consistency, its advantage
lies in no need to design an ad-hoc loss for each specific data
generation process, which is generally required in the consistent
methods. In conclusion, robustness is a common and critical
theoretical guarantee in supervised learning, but the mechanism
by which it might be achieved remains barely understood in PLL.
To the best of our knowledge, this is the first work to analyze
the robustness of PLL.

IV. METHODOLOGY

In this section, we propose a family of APL loss functions for
PLs, and introduce the generation processes of PLs.

A. A Family of Average PL (APL) Losses

In this paper, we propose a family of loss functions named the
average PL (APL) losses following the principled ABS:

$$\tilde{\ell}(f(x), s) = \frac{1}{|s|} \sum_{i \in s} \ell(f(x), i),$$
where $| \cdot |$ represents the cardinality. Our learning formulation
is built on a simple scheme that combines multiple multi-class
losses on the individual candidate. For example, we can use the
GCE or CCE loss as the component $\ell$. If $s$ is a singleton, the
APL loss reduces to the ordinary multi-class loss. The idea of
the APL losses comes from a practically motivated process proposed
by Feng et al. [18]: they assumed that a candidate-label set is
feature-independent and uniformly drawn given a specific true
label, i.e., $p(s|y, x) = p(s|y) = \text{const.}$, if $y \in s$, and $p(s|y) = 0$
otherwise. The generation process of noise-free PLs can thus be
formalized as $p(s|x) = \sum_y p(s|y)p(y|x) \propto \sum_{y \notin s} p(y|x)$.

Then we could consider replacing the posterior with a loss and

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obtain $\hat{\ell}(f(x), s) \propto \sum_{y \in S} \ell(f(x), y)$. It inspires the formula of the APL losses, and the normalization term $1/|s|$ breaks the bias to training data with more candidate labels. The APL losses encourage the larger outputs on candidate labels, while do not explicitly guarantee the true label has the biggest score. Ideally, a “nice” loss $\ell$ can drive up the output of the true label implicitly resorting to the inductive bias, while a “bad” loss results in an inability to disambiguate. Thus, the issue now is that which multi-class loss functions can bound $\mathcal{R}(f^*) - \mathcal{R}^*$ (or $\overline{\mathcal{R}}(f^*) - \overline{\mathcal{R}}^*$), that is, make our ABS method with the APL loss PL-robust.

Let us give an motivating example. $\{z_1, z_2\} \in S$ are two candidate labels of an instance $x$, and $z_1$ is true. Then the APL loss of $f$ on this sample is $\hat{\ell}(f(x), s) = \frac{1}{2}[\ell(f(x), z_1) + \ell(f(x), z_2)]$. We would like to increase $g_{z_1}(x)$ to get it close to 1 so that $g_{z_2}(x)$ is decreased, signifying $f$ successfully remembers the true label without the interference of $z_2$. Paradoxically, because all candidate labels contribute to minimizing $\hat{\ell}$, neither $\ell(f(x), z_1)$ nor $\ell(f(x), z_2)$ should be too large. Intuitively, if $\ell$ has an upper bound, then the value of $\ell$ is acceptable even if $g_{z_2}(x)$ is close to 0. But if this is not the case, the optimization algorithm must keep $g_{z_2}(x)$ not too small to ensure that $\hat{\ell}(f(x), z_2)$ is not too small, then memory for the true labels is hindered. The empirical observations in Section I also confirm this inference.

We investigate a series of non-negative multi-class loss functions and prove that the APL losses with bounded multi-class loss functions$^1$ are robust to (both noise-free and noisy) PL data in Section V.

**Definition 3.** We say a multi-class loss function is bounded if for any classifier $f$ and input $x \in \mathcal{X}$, the sum of losses over all classes is bounded by certain constants $C_1$ and $C_2$:

$$
C_1 \leq \sum_{i=1}^{k} \ell(f(x), i) \leq C_2. \quad (3)
$$

Especially, if $C_1 = C_2 = C$, i.e., $\sum_{i=1}^{k} \ell(f(x), i) = C$, the loss function is said to be symmetric.

We should point out that the loss of a bounded loss function on any class is also bounded: $0 \leq \ell(f(x), i) \leq U$, $\forall i \in \mathcal{Y}$, where $U$ is a constant. We examine widely-used loss functions and list their bounds in Table I. We use a one-hot representation for each label, i.e., if the label $y = i$, its label vector is represented as $e_i$, where the $j$th element is given by $e^j_i = 1$ if $i = j$, otherwise 0. Then for symmetric losses, $\ell(f(x), j) = C/(k-1), \forall j \neq i$.

### B. Data Generation Processes

To provide the main insights, we have to propose some assumptions on the data generation processes. In the following, we formally establish two general problem settings for the generation processes of noise-free PLs and noisy PLs, and for each of them we further propose several particular implementations motivated by realistic scenarios. We follow the assumption in prior reliable PLL works $^{[18],[36],[61]}$ and the classical feature-independent model of label noise $^{[24],[45],[50],[53]}$ that the observation is conditionally independent of input given the true label, as a result there are $s \perp x | y, p(s|x, y) = p(s|y)$. Then the density of corrupted distribution is formulated as $p(x, s) = \sum_{y \in \mathcal{Y}} p(s|y)p(x, y)$.

#### 1) Filtered Sampling Process for Noise-Free PLs: In a pioneering study involving PLL generation processes, the PL is assumed to be independently and uniformly sampled given a specific true label $^{[18]}$. It is inspired by a real-world cost-saving application of labeling: without any prior knowledge, the labeling system generates a random PL for each sample and asks the human annotators whether the set contains the true label. Resampling PLs for samples for which the annotators answered “NO”. While it would be easy to make the labeling system have some rudimentary knowledge, for example, “salmon” and “spacecraft” do not usually appear in the same set. Then we generalize the uniformly sampling assumption and propose the Filtered sampling process. Formally speaking, given a specific true label, a noise-free PL is assumed to be sampled as a whole:

$$
p(s|y) = \begin{cases} 
\eta_y^k & \text{if } y \in S, \\
0 & \text{if } y \notin S,
\end{cases} \quad (4)
$$

where $0 \leq \eta_y^k \leq 1$ is the sampling probability of the label set $s$ given the true label $y$, and $\sum_{y \in \mathcal{Y}} \eta_y^k = 1$.

This generation process can be written in the form of a transition matrix $^{[24]}$. We enumerate all the label sets $s$ in the PL space $\mathcal{S}$ and specify an index $l$ for each set, i.e., $l_s \in \mathcal{S}$ $(i \in [2^k - 2])$. By this notion, we summarize all the probabilities into a PL transition matrix $Q^{k \times (2^k - 2)}$, where $Q_{lij} = p(s = l_j|y = i)$. Further taking into account the assumed data distribution in (4), we can instantiate $Q$ as $Q_{ij} = \eta_y^j$ if $i \neq l_j$, otherwise $Q_{ij} = 0$. Then for all $j \in [2^k - 2]$, there is $p(s = l_j|x) = \sum_{i \in l_j} p(s = l_j|y = i)p(y = i|x)$. Thus we have

$$
p(x, s) = Q^\top p(x, y), \quad (5)
$$

where $^\top$ denotes the transpose. **Flipping Model.** In the real world, however, there are complex and varying correlations between categories, making it common to get some similar categories mixed up, so that some combinations of labels appear more frequently than others. The probability of an incorrect label appearing in the PL of a sample depends on how similar it is to its true label. We therefore propose the Flipping model as a means of representing this more realistic situation, where a noise-free PL is supposed to be generated by adding each label to the candidate-label set independently:

$$
p(s|y) = M \prod_{i \in S} \eta_i^y \prod_{i \notin s} (1 - \eta_i^y), \quad (6)
$$

where

$$
\eta_i^y = p(i \in s|y), \forall i \in \mathcal{Y}, \quad M = 1/ \left(1 - \prod_{i \neq y} \eta_i^y\right).
$$

$\eta_i^y$ is the flipping probability that depicts the probability of $i$-label being included into the candidate-label set given the specific class label $y$, and $\eta_y^0 = 1$. For $i \neq y$, it satisfies $0 \leq \eta_i^y < 1$. $M$ excludes the set whose cardinality equals $k$ by re-sampling.
Similarly, the PL transition matrix can be formulated as $Q^{k \times k}$ where $Q_{ij} = p(j \mid s \mid y = i) = \eta_{y}^{i}$ and the diagonal elements of $Q$ are all 1. If $q(x)$ is a $k$-dimension vector where the $j$th element $q_{j}(x)$ is the probability $p(j \mid s \mid x)$, then

$q(x) = Q^{T} p(y \mid x),$

$p(x, s) = M \prod_{i \in S} q_{i}(x) \prod_{j \notin s} (1 - q_{j}(x)) p(x).$ (7)

The Flipping model is a well-established approach for characterizing the data generation process in weakly supervised learning problems represented by NLL. This model emphasizes the commonalities among categories and quantifies the likelihood that an error category is indistinguishable from the true category, parameterized by a class-level transition matrix. In contrast, in PLL, the Filtered sampling process serves as the “relabel complete” Flipping model, as annotations in PLL form a transition matrix at the collection (i.e., PL) level. However, the Filtered sampling process typically fails to decouple the flipping probabilities of individual labels, making it challenging to discern similarities among categories. Consequently, special consideration of the Flipping model remains necessary.

2) Global Sampling Process for Noisy PLs: As a reminder, the sampling process entails manual filtering of non-conforming label sets and subsequent resampling, leaving room for potential noisy PLs in the event of errors by human annotators. Consequently, all elements within the PL space possess a non-zero probability of being sampled:

$p(s \mid y) = \eta_{y}^{s}, \forall s \in S,$ (8)

where $0 \leq \eta_{y}^{s} \leq 1$. If the sampling probability for all noise-free PLs is $(1 - \gamma)/(2^{k-1} - 1)$, and that for all noisy PLs equals $\gamma/(2^{k-1} - 1)$, then the sampling probability under the Global sampling process is said to be uniform. In addition, the density $p(x, s)$ takes the same form as (5) while even if $i \notin s$, $Q_{ij}$ may be larger than 0.

Next we discuss two specific generation patterns of the Global sampling process in terms of how noise-free PLs are contaminated, i.e., the true class is obfuscated by another similar class, or noise-free PL is deliberately corrupted, respectively.

Confusing Model. In this type of setting, the true class was (accidentally) confused with other (similar) classes, leading to the misuse of an incorrect label as the original true label in the PL generation process. Therefore, the Confusing model consists of two steps.

First, the true label is corrupted. Suppose the class-conditional noise (CCN) model [24, 45, 50, 53, 64]—the most widely-used model for noisy label classification—is applied, where each instance from class $y$ has a fixed probability of being assigned to label $i$, that is

\[
\tilde{y} = \begin{cases} y & \text{with probability } 1 - \gamma_{y}, \\ i, i \in \mathcal{Y}, i \neq y & \text{with probability } \gamma_{y}^{i} \end{cases}
\]

where $0 \leq \gamma_{y} \leq 1$ is the label noise rate. The noise is said to be uniform if $\gamma_{y}^{i} = \gamma$ and $\gamma_{y}^{i} = \gamma/(k - 1)$, otherwise it is said to be asymmetric. The corrupting step can be formalized by the noise transition matrix $T$ [53], where $T_{ij} = p(\tilde{y} = j \mid y = i)$. Second, the corrupted label $\tilde{y}$ serves as the true label $y$ to generate candidate labels, which signifies we require noisy PLs to contain $\tilde{y}$ in the same way that noise-free PLs must contain $y$. Accordingly, the following equation holds:

\[
p(x, s) = \sum_{y, \tilde{y} \in \mathcal{Y}} p(s \mid \tilde{y}) p(\tilde{y} \mid y) p(y \mid x) p(x) = \sum_{j=1}^{k} \sum_{i=1}^{k} \mathcal{T}_{ij}^{T} p(y = i \mid x) p(x),
\]

where $p(s \mid y)$ can be expanded into the form of (4) or (6).

Destructing Model. We believe that noise-free PLs can also be (intentionally) destructed. For example, there may exist spammers who deliberately choose label sets that are totally irrelevant to the tasks. Hence we propose the Destructing model that also contains two steps.

Noise-free PLs are first generated by the Filtered sampling process, followed by taking its complement with the set flipping

| Loss function | Bound of loss | Bound of the sum of losses over all classes |
|---------------|--------------|-------------------------------------------|
| MAE $\ell(f(x), i) = |e_i - f(x)|| 0 \leq \ell(f(x), i) \leq 2 | \sum_{i=1}^{k} \ell(f(x), i) = 2k - 2$
| MSE $\ell(f(x), i) = |e_i - f(x)|| 0 \leq \ell(f(x), i) \leq 2 | k - 1 \leq \sum_{i=1}^{k} \ell(f(x), i) \leq 2k - 2$
| RCE $\ell(f(x), i) = - \sum_{i=1}^{k} g_i(x) \log e_i | 0 \leq \ell(f(x), i) \leq -A, A < 0 | \sum_{i=1}^{k} \ell(f(x), i) = A - Ak$
| GCE $\ell(f(x), i) = \frac{1}{\log q_i(x)} | 0 \leq \ell(f(x), i) \leq \frac{1}{q} \text{ with } q \in [0, 1] | k \log k \leq \frac{1}{\gamma} \sum_{i=1}^{k} \ell(f(x), i) \leq k(1)(k - 1)(k - 2) / (k - 1)^2$
| PCE $\ell(f(x), i) = \left\{ \begin{array}{ll} -\gamma g_i(x) + \log \tau + 1, & \text{if } g_i(x) \leq \frac{1}{\gamma} \\ -\log g_i(x), & \text{otherwise} \end{array} \right. | 0 \leq \ell(f(x), i) \leq \log \tau + 1, \tau > 1 | k\gamma = \frac{1}{\gamma} \sum_{i=1}^{k} \ell(f(x), i) \leq (k - 1)(k - 1)(k - 1)(k - 2) / (k - 1)^2$
| CCE $\ell(f(x), i) = \log q_i(x) | \ell(f(x), i) \geq 0 | \ell(f(x), i) > 0, \tau > 0 | \text{Unbounded}$
| PL $\ell(f(x), i) = -(1 - g_i(x)) \log q_i(x) | \ell(f(x), i) > 0, \tau > 0 | \text{Unbounded}$

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rate $\gamma_s$: 
\[
\gamma_s = \begin{cases} 
  s & \text{with probability } 1 - \gamma_s, \\
  \bar{s} & \text{with probability } \gamma_s, 
\end{cases} 
\] 
where $0 \leq \gamma_s \leq 1$. Constructing the complement transition matrix $T^* = (2^k - 1) \times (2^k - 1)$, each row (column) of which represents a fixed label set in the PL space. Let $T_{ij} = \gamma_s$ if the $i$th and the $j$th label set are complementary to each other, $T_{ij} = 1 - \gamma_s$ for all $i \in [2^k - 2]$, and 0 otherwise. Then the density function of the PL distribution $p(x, s)$ is multiplied by $T$ on the basis of (5) or (7).

Similar to the noise-free scenario, the Global sampling process provides a thorough description of the sampling possibilities for all PLs, whereas the Confusing and Destructing models, which are particular instances of the Global sampling process, focus primarily on elucidating the techniques and approaches used to correct true labels and generate PLs.

V. THEORETICAL RESULTS

The studied generation processes of PLs allow us to theoretically understand the properties of the APL losses. In this section, we detail sufficient conditions under which multi-class loss functions make PLL with the APL losses robust in various scenarios, and conclude some instructive findings.

A. Robustness to Noise-Free PLs

Lemma 1. Any symmetric loss satisfies $\hat{f}^* = \hat{f}^*$, and any bounded loss satisfies $0 \leq \bar{R}(\hat{f}^*) - \bar{R}^* \leq \frac{A(C_2 - C_1)}{4A}$, where 
\[
A = \frac{1}{2^m - 1} \sum_{j=1}^{m-1} \frac{1}{2^j - 1} \sum_{i=1, i \neq y} \eta_y^i.
\]
We introduce two quantities, $A$ is the expectation of a weighted sum of $p(y \in s')$ over $p(x, y)$, and $A'$ is the weighted sum of $p(i \in s, i \neq y)$. Theorem 1 establishes that under certain conditions, the risk of the Bayes classifier on PL data approaches the minimal PLL risk w.r.t. bounded losses. The tighter the bound of the bounded loss, the more robust the APL loss, and the extreme case is achieved by the symmetric loss which leads to statistical consistency (refer to Section III-B). The crucial condition for PL-robustness concerns the sampling probability. Since $\sum_{y \in s} \eta_y^i = 1$, the condition $\sum_{s \in I} \eta_s^i < 1, \forall i \neq y$ signifies that any label other than the true one may not necessarily included in the candidate-label set, i.e., the domination of true labels. Another constraint, $\bar{R} = 0$, implies that classes are separable in fully-supervised classification if the multi-class loss $\ell$ is classification-calibrated. Note that experimental results later demonstrate that even if this constraint is not satisfied, bounded losses still exhibit good empirical PL-robustness.

In Lemma 1, the uniform sampling probability has already ensured the domination of true labels, and the separability constraint $\bar{R} = 0$ is eliminated, which implies that even in the stochastic scenario, learning with the APL losses can be PL-robust. Therefore, in this case, the robustness is satisfied without any constraint. In addition, the excess 0-1 risk is bounded, which signifies that the optimal PLL classifier approaches the Bayes classifier, so that the PL-robustness is better guaranteed.

Moreover, if we take into account the shared information among categories the utilize the Flipping model to formalize the generation process, we can obtain the subsequent PL-robustness conditions.

Corollary 2. With the Flipping model, suppose $\bar{R}^* = 0$, then
1) for any symmetric loss, $\hat{f}^* = f^*$;
2) for any bounded loss, $0 \leq \bar{R}(\hat{f}^*) - \bar{R}^* \leq MA(C_2 - C_1)$, where 
\[
A = \frac{1}{2^m - 1} \sum_{j=1}^{m-1} \frac{1}{2^j - 1} \sum_{i=1, i \neq y} \eta_y^i 
\]
and $A'$ is the weighted sum of $p(i \in s, i \neq y)/p(x, y)$ under the Global sampling process with uniform sampling probability.

Theorem 1. With the Filtered sampling process, suppose $\bar{R}^* = 0$ and $\forall i \neq y, \sum_{s \in I} \eta_s^i < 1$, then
1) for any symmetric loss, $\hat{f}^* = f^*$;
2) for any bounded loss, $0 \leq \bar{R}(\hat{f}^*) - \bar{R}^* \leq MA(C_2 - C_1)$, where 
\[
A = \frac{1}{2^m - 1} \sum_{j=1}^{m-1} \frac{1}{2^j - 1} \sum_{i=1, i \neq y} \eta_y^i 
\]
and $A'$ is the weighted sum of $p(i \in s, i \neq y)/p(x, y)$ under the Global sampling process with uniform sampling probability.

Comparing this corollary versus Theorem 1, we observe that their sufficient conditions and upper bounds on the risk difference differ in notation, but are conceptually similar. Notably, the flipping probability is not constrained due to the difference differ in notation, but are conceptually similar. No-

B. Robustness to Noisy PLs

In this section, we present formal statements for the Global sampling process, and provide intuitive statements for the specific models of noisy PLs, as articulated in Corollary 4, with more formal mathematical details deferred to Appendix A, available online.

Lemma 2. Any symmetric loss satisfies $\hat{f}^* = \hat{f}^*$, and any bounded loss satisfies $0 \leq \bar{R}(\hat{f}^*) - \bar{R}^* \leq \frac{A(C_2 - C_1)}{4A}$, where 
\[
A = \frac{1}{2^m - 1} \sum_{j=1}^{m-1} \frac{1}{2^j - 1} \sum_{i=1, i \neq y} \eta_y^i 
\]
and $A'$ is the weighted sum of $p(i \in s, i \neq y)/p(x, y)$ under the Global sampling process with uniform sampling probability, if label noise rate $\gamma < 1/2$.

Theorem 3. With the Global sampling process, suppose $\bar{R}^* = 0$ and the domination relations hold: $d(y) > d(i)$ whenever $d(i) \neq d(y)$, where $d(i)$ is defined as 
\[
d(i) = \frac{1}{2^m - 1} \sum_{j=1}^{m-1} \frac{1}{2^j - 1} \sum_{i=1, i \neq y} \eta_y^i 
\]
and $A'$ is the weighted sum of $p(i \in s, i \neq y)/p(x, y)$ under the Global sampling process with uniform sampling probability, if label noise rate $\gamma < 1/2$.

Theorem 3. With the Global sampling process, suppose $\bar{R}^* = 0$ and the domination relations hold: $d(y) > d(i)$ whenever $d(i) \neq d(y)$, where $d(i)$ is defined as 
\[
d(i) = \frac{1}{2^m - 1} \sum_{j=1}^{m-1} \frac{1}{2^j - 1} \sum_{i=1, i \neq y} \eta_y^i 
\]
and $A'$ is the weighted sum of $p(i \in s, i \neq y)/p(x, y)$ under the Global sampling process with uniform sampling probability, if label noise rate $\gamma < 1/2$. The condition for robustness to noisy PLs tightens the constraint on the dominance relationship compared to noise-free PLs, from the sum of probabilities to the weighted sum of probabilities. Whereas, with the uniform sampling probability, the constraint is relatively loose: the probability of outputting noise-free PLs
exceeds fifty percent, requiring only slightly more domain knowledge than a completely random labeling system.

To encompass a variety of cases, we summarize the Confusing and Destructing models in the following corollary and then discuss each in detail.

**Corollary 4.** With the Confusing model or the Destructing model, suppose \( R^* = 0 \) and the domination relations hold: the expectation of the weighted sum of \( p(y \in \mathcal{s}|y) \) is always larger than that of \( p(i \in \mathcal{s}|y), \forall i \neq y \) over \( p(x|y), \) then

1) for any symmetric loss, \( f^* = f; \)
2) for any bounded loss, \( 0 \leq R(f^*) - R^* \leq A(C_2 - C_1), \)

where \( A \) is a constant associated with \( p(y \in \mathcal{s}|x). \)

For the Confusing model, we assume uniform cases in the generation of candidate labels to simplify the process. This simplification degenerates the domination relations to \( \gamma_{iy} < 1 - \gamma_y, \forall i \neq y, \) which we were a little surprised to find that it becomes identical to the condition for asymmetric-noise-tolerance (Theorem 3 in [20]). Further supposing that the true labels are corrupted uniformly, we have \( \gamma < (k - 1)/k. \) This is the same as the noise-tolerant condition under uniform noise (Theorem 1 in [20]). These findings indicate that the robustness condition to noisy PLs is exclusively determined by the label noise rate as long as the candidate labels are generated uniformly. Moreover, the constraint \( R^* = 0 \) is removed and \( R(f^*) - R^* \) is bounded in this case.

Next we probe into the Destructing model through similar renderings. Assuming uniform generation of candidate labels, the domination relations are reduced to \( \gamma_s < 1/k, \forall s. \) If the probability of destructing every candidate-label set is also uniform, the set flipping rate is also the unreliability rate, namely \( \gamma_s = \gamma, \forall s. \) Then we again show that the PL-robustness condition is solely dependent on the level of unreliability in the candidate-label set.

**Remarks.** Our theoretical analysis on the APL losses for learning with PL data reveals several critical conditions for achieving PL-robustness. We now make several observations about all the theorems:

- The key factor that makes the APL losses PL-robust is consistent across all scenarios: the weighted sum of the probabilities of the true labels associated with the instances dominates;
- For noisy PL data, if the candidate labels are generated uniformly, the robustness condition is exclusively determined by the degree of unreliability rate;
- In the uniform case, the excess 0-1 risk w.r.t. the clean distribution of the optimal PLL classifier can be upper bounded in proportion to the difference between the upper and lower bounds of the loss function \( C_2 - C_1. \) A smaller difference implies a smaller upper bound on the excess risk, implying a more robust APL loss, and then the most desirable situation (statistical consistency) can be achieved by the symmetric losses. Its constant of proportionality increases with the probability of incorrect labels, accord with the intuition that higher stochasticity makes learning difficult;
- In the more general cases, when learning with a bounded loss, we can only bound the excess \( \ell \)-risk under the corrupted distribution, whose upper bound, up to a constant less than 1, equals to \( C_2 - C_1. \)

The above theoretical findings provide guidance to the design of losses of ABS. Note that IBS is heuristic and not really ERM-based (refer to Section. VI-C), and on this account, the robustness of IBS could hardly be proven.

**C. Estimation Error Bound**

Let us review the relation between robustness and consistency again. Now we have proved the condition that bounds RHS2 in (1) (assuming \( R^* \) is instantiated to be a DNN). Then we establish the estimation error bound and show that as the number of training data approaches infinity, RHS1 is also bounded.

Suppose \( \mathcal{G}_y \) be a class of real functions, and \( \mathcal{F} = \oplus_{y \in [k]} \mathcal{G}_y \) be a \( k \)-valued function class. Assume there are \( C_f > 0 \) and \( C_\ell > 0 \) such that \( \sup_{f \in \mathcal{F}} \|f\|_\infty \leq C_f \) and \( \sup_{x \in \mathcal{X}: f \in \mathcal{F}, y \in [k]} \ell(f(x), y) \leq C_\ell, \) and assume \( \ell(f(x), y) \) is \( \rho \)-Lipschitz continuous for all \( \|\|_\infty \leq C_f. \) The Rademacher complexity of \( \mathcal{G}_y \) over \( p(x) \) with sample size \( n \) is defined as \( \mathfrak{R}_n(\mathcal{G}_y) \) [5, 48]. Then we have the following estimation error bound.

**Theorem 5.** For any \( \delta > 0, \) we have with probability at least \( 1 - \delta, \)

\[
\hat{R}(f) - \hat{R}(f^*) \leq 4\sqrt{2}k \mathfrak{R}_n(\mathcal{G}_y) + 2C_\ell \sqrt{\frac{\log(2/\delta)}{2n}}.
\]

As \( n \to \infty, \) \( \mathfrak{R}_n(\mathcal{G}_y) \to 0 \) for all parametric models with a bounded norm such as DNNs trained with weight decay [39], which signifies \( \hat{R}(f) \to \hat{R}(f^*). \)

**VI. EXPERIMENTAL FINDINGS**

In this section, we provide some empirical understandings of our ABS method, experimentally validating our theoretical findings on benchmark datasets, which then enlighten an improvement of IBS methods. The implementation is based on PyTorch [52] and experiments were carried out with NVIDIA Tesla V100 GPU.

**A. Empirical Understanding of APL Losses**

**Our ABS Method With Bounded Loss Functions Are Robust.** We first run a set of experiments on MNIST [34] and CIFAR-10 [32] to verify whether our ABS method with bounded losses is robust to both noise-free PLs and noisy PLs, whereas they are not with unbounded losses. We generate noise-free PLs by the Flipping model with a uniform flipping probability of 0.1, and then noisy PLs by the Confusing model with \( \gamma = 0.3. \) On each dataset, we train two networks using the APL losses with different multi-class losses, e.g., bounded versus unbounded ones in Table I. The RCE loss is, up to a constant of proportionality, equivalent to the MAE loss and omitted. We set the focusing parameter 0.5 for the FL loss. Detailed settings are in Section VI-B.

The test accuracy with different losses is presented in Fig. 2. As we have theoretically proved, learning with bounded losses is robust: after reaching a peak in test accuracy, their test accuracy is relatively flat throughout the training process. However,
unbounded losses exhibit significant overfitting in most cases. Specifically, the symmetric loss MAE has the smoothest curve, while other bounded losses would be slightly overfitting in difficult learning scenarios. The same results are shown across different datasets, under different data settings, with different models. In general, the more difficult the learning scenario is (e.g., harder datasets and weaker supervised information), the larger the gap between bounded loss and unbounded loss is, because unbounded losses overfit more severely.

How do the Models Fit the Candidate Labels When Training With the APL Losses? As we discussed before, ABS methods are free from identifying the true labels during training. One may wonder how ABS methods learn from PL data. This raises the question: is the learned model able to identify the true label of the training samples? We investigate this problem by looking at the confidence margin between the model’s output on the true label and the maximum output on the other labels, i.e., confidence margin($x_i$) = $g_{y_i}(x_i) - \max_{j \neq y_i} g_j(x_i)$. The larger the margin is, the greater the likelihood that the model will successfully identify the true label for the training samples is. In Fig. 3, we illustrate the mean confidence margin over the training set. We find that the margins trained with bounded losses are generally much higher than those trained with unbounded losses. This means that although our ABS method does not explicitly disambiguate the candidate-label sets during the training phase, our ABS method with bounded losses is still able to robustly fit the true labels against the interference of other candidates, thereby explaining the good prediction performance in the test set.

Fig. 2. Test accuracy of our ABS method with bounded versus unbounded losses on benchmarks.

Fig. 3. Confidence margin of our ABS method with bounded versus unbounded losses on benchmarks.
TABLE II
MEANS ± STANDARD DEVIATIONS OF TEST ACCURACY IN PERCENTAGE WITH DIFFERENT DATA GENERATIONS

| Dataset       | Model         | Case | Bounded | Unbounded |
|---------------|---------------|------|---------|-----------|
|               | MEANS ± MSE  | GCE  | PCE     | CCE       | FL        |
| MNIST         | Linear        |      |         |           |           |
|               | 1 91.17 ± 0.07| 87.61±0.07 | 90.40±0.12 | 86.28±0.40 | 85.13±0.29 | 85.10±0.24 |
|               | 2 92.25±0.08  | 82.03±0.10 | 93.33±0.06 | 92.71±0.08 | 89.51±0.17 | 89.71±0.13 |
|               | 3 91.90±0.13  | 89.90±0.10 | 91.86±0.10 | 91.25±0.19 | 87.28±0.27 | 87.08±0.27 |
|               | 4 90.37±0.19  | 82.70±0.49 | 85.87±0.48 | 80.61±0.41 | 67.78±1.19 | 67.63±0.84 |
| MLP-5         | Linear        |      |         |           |           |
|               | 1 94.44 ± 0.19| 82.28±0.83 | 92.77±0.21 | 87.62±0.27 | 84.80±0.50 | 84.15±0.73 |
|               | 2 96.71 ± 0.05| 95.16±0.19 | 96.49±0.15 | 95.92±0.13 | 93.08±0.30 | 92.62±0.18 |
|               | 3 95.87 ± 0.85| 80.71±0.52 | 95.34±0.19 | 91.14±0.25 | 77.68±0.53 | 75.66±0.74 |
|               | 4 93.18 ± 0.47| 89.12±1.03 | 89.19±0.47 | 86.52±0.82 | 77.57±1.32 | 76.45±0.91 |
| Fashion-MNIST | MLP           |      |         |           |           |
|               | 1 86.37 ± 0.64| 81.26±0.47 | 84.29±0.10 | 76.01±0.49 | 51.29±0.83 | 49.36±0.48 |
|               | 2 87.52 ± 0.09| 84.49±0.27 | 87.32±0.31 | 85.47±0.05 | 74.10±0.64 | 73.50±0.24 |
|               | 3 85.74 ± 0.27| 81.62±0.33 | 84.42±0.37 | 80.33±0.51 | 53.09±0.74 | 52.75±0.81 |
|               | 4 82.41 ± 0.25| 73.49±0.53 | 66.83±0.87 | 67.25±0.87 | 25.76±0.38 | 22.01±0.53 |
| Kuzushiji-MNIST | MLP         |      |         |           |           |
|               | 1 63.14±0.14  | 59.65±0.62 | 65.58±0.32 | 58.78±0.46 | 54.90±0.61 | 55.02±0.48 |
|               | 2 64.35±0.09  | 68.54±0.31 | 65.04±0.20 | 67.40±1.93 | 63.79±0.34 | 63.74±0.44 |
|               | 3 63.55±0.22  | 65.26±0.21 | 64.24±0.29 | 67.75±0.17 | 59.82±0.36 | 60.16±0.73 |
|               | 4 60.34±0.67  | 56.16±0.80 | 51.98±0.44 | 51.47±0.80 | 36.48±0.59 | 35.76±0.31 |
| CIFAR-10      | MLP           |      |         |           |           |
|               | 1 81.72 ± 0.50| 64.60±1.06 | 75.71±0.15 | 57.17±0.22 | 36.07±0.16 | 26.13±0.46 |
|               | 2 86.99 ± 0.25| 74.24±0.13 | 86.59±0.24 | 80.90±0.41 | 66.06±0.51 | 64.81±0.41 |
|               | 3 78.66±0.25  | 69.90±0.61 | 79.26±0.54 | 72.89±0.18 | 56.43±0.88 | 54.96±0.64 |
|               | 4 66.61 ± 0.29| 54.93±0.18 | 43.06±0.44 | 42.58±0.43 | 23.93±0.44 | 23.35±0.63 |
|               | ConvNet       |      |         |           |           |
|               | 1 41.46 ± 0.72| 33.11±0.69 | 31.98±0.40 | 31.10±0.42 | 22.47±0.49 | 22.75±0.28 |
|               | 2 48.69 ± 0.25| 38.99±0.42 | 46.00±0.30 | 42.94±0.44 | 36.81±0.22 | 36.37±0.35 |
|               | 3 40.83 ± 0.48| 29.05±0.49 | 35.30±0.30 | 30.90±0.49 | 28.04±0.34 | 27.31±0.18 |
|               | 4 31.62 ± 0.60| 25.52±0.55 | 20.75±0.35 | 21.10±0.37 | 13.51±0.34 | 13.70±0.35 |

B. Evaluation on Benchmark Datasets

Setup. Experiments were conducted on four widely-used benchmark datasets including MNIST, Fashion-MNIST [65], Kuzushiji-MNIST [11], and CIFAR-10. On each dataset, we generated PLs by (Case 1) the Filtered sampling process with a uniform sampling probability; (Case 2) the Flipping model with a uniform flipping probability of 0.1; (Case 3) the Confusing model where the label noise rate equals 0.3 and the candidate labels were generated according to Case 2; (Case 4) the Destructing model where the candidate labels were generated according to Case 1 and the set flipping rate equals 0.3. We leave out 10% of the corrupted training samples as a validation set, which is for model selection. Further experiments on the generation processes with different parameters can be found in Appendix C, available online.

We employed various base models including a linear-in-input model (Linear), a 5-layer perceptron (MLP), and a 12-layer convolutional neural network (ConvNet) [25]. Linear was trained on
TABLE III
PRODEN & OUR ABS METHOD ON CIFAR DATASETS

| Dataset   | Model  | Case | E: | A: | X | √ | X | √ | X | √ | Our ABS |
|-----------|--------|------|----|----|---|---|---|---|---|---|---------|
| CIFAR-10  | MLP    | 1    | A: |    | 48.28 ± 0.62 | 48.33 ± 0.54 | 48.77 ± 0.26 | 48.54 ± 0.36 | — |         |
|           |        | 2    |    |    | 51.49 ± 0.30 | 52.17 ± 0.21 | 52.31 ± 0.32 | 52.70 ± 0.14 | — |         |
|           |        | 3    |    |    | 38.52 ± 0.16 | 44.27 ± 0.13 | 39.00 ± 0.38 | 46.67 ± 0.30 | 45.86±0.14 |         |
|           |        | 4    |    |    | 29.19 ± 0.63 | 34.04 ± 0.56 | 29.30 ± 0.43 | 35.12 ± 0.42 | 34.54±0.36 |         |
| CIFAR-100 | ConvNet| 1    |    |    | 85.38 ± 0.11 | 85.52 ± 0.25 | 85.93 ± 0.31 | 86.09 ± 0.20 | — |         |
|           |        | 2    |    |    | 88.62 ± 0.19 | 88.29 ± 0.33 | 89.42 ± 0.17 | 89.05 ± 0.29 | — |         |
|           |        | 3    |    |    | 64.59 ± 0.70 | 73.05 ± 0.31 | 66.86 ± 0.72 | 75.99 ± 0.39 | 76.72±0.12 |         |
|           |        | 4    |    |    | 47.35 ± 0.47 | 53.47 ± 0.38 | 50.77 ± 0.82 | 56.68 ± 0.49 | 55.50±0.24 |         |
| CIFAR-100 | ConvNet| 2    |    |    | 54.49 ± 0.46 | 59.90 ± 0.53 | 58.31 ± 0.32 | 60.60 ± 0.22 | — |         |
|           |        | 3    |    |    | 34.78 ± 0.31 | 42.04 ± 0.22 | 36.27 ± 0.43 | 43.67 ± 0.22 | 44.75±0.15 |         |
|           |        | 4    |    |    | 37.37 ± 0.32 | 47.29 ± 0.31 | 39.55 ± 0.24 | 47.86 ± 0.20 | 46.13±0.24 |         |

E and A stand for early stopping and robust warm start, respectively. The best and equivalent based on the paired t-test at the significance level 5% are shown in boldface by comparing the 1st and 3rd columns, the 2nd and 4th columns. “—” means that we skipped the experiments under the reliable PLL setting. The best combination is underline.

MNIST-like datasets, ConvNet was trained on CIFAR-10, and MLP was trained on all datasets. The optimizer was stochastic gradient descent with momentum 0.9. We trained each model 500 epochs with the mini-batch size set to 256, and recorded the test accuracy of the hyper-parameters (learning rate and weight decay) with the best validation accuracy. We did not use any manual learning rate decay and early stopping. Results. Tables II shows the test accuracy over 5 trials. The best and comparable methods based on the paired t-test at the significance level 5% were highlighted in boldface. We can see that the bounded loss always outperforms the unbounded loss, especially on the complex models. In difficult scenarios, i.e., unreliable PLL, the accuracy of the complex models trained with bounded losses is almost always better than their linear counterpart, but unbounded losses can make the complex models overfit very badly on some tasks, causing their performance to become worse.

C. Enhancing IBS Methods With ABS

We revisit the SOTA IBS methods PRODEN [41], RC [18], and LW [61]. Their typical learning objective is as follows:

\[
\mathcal{R}(f; \ell) = \frac{1}{n} \sum_{i=1}^{n} \left[ \sum_{j \in s_i} w_j \ell(g_j(x_i), j) + \sum_{j \notin s_i} w_j \ell(g_j(x_i), j) \right],
\]

where \( w_j \) is a weight for \( j \in [k] \). The weights of the labels that are more likely to be true are progressively increased. Generally speaking, they initialize the weights to be \( w_j = 1/|s_i| \), \( \forall j \in s_i \).

They train a learning model \( f \) with the uniform weights for several epochs for a warm start, and then update \( w \) and \( f \) seamlessly for the remaining epochs. We highlight that uniform weights are necessary to break the circular dependency existing between \( w \) and \( f \): \( f \) needs to be trained with reasonable \( w \) and \( w \) needs to be estimated by well-trained \( f \). The success of the algorithms is built on the observation that even if each sample has multiple candidate labels, \( f \) will remember the true one first [3]. Thus they adjust \( w \) by the output of \( f \). It indicates that an IBS method has to be pretrained a little in an ABS manner.

While we note that they always use the CCE loss in both the ABS-style phase and the subsequent IBS-style phase, which could potentially select incorrect true labels at the very beginning and negatively affect the model training. Therefore, we introduce an enhanced principle to incorporate our theoretical findings into existing IBS methods: training the learning model with a robust warm start to avoid overfitting. We replace their loss functions of the first 20 epochs with the APL loss with MAE, and then switch back to their original objective function. The hyper-parameters are tuned according to the original methods.

We considered Case 1, 2, 3, and 4 for CIFAR-10, and Case 3, 2, and 4 where the candidate labels were generated by the Flipping model with a uniform flipping probability of 0.01 for CIFAR-100 [32]. We used the same training/validation setting, models, and optimizer as in Section VI-B. We summarize the results without/with early stopping of PRODEN in Table III, which means that we report the last epoch or the epoch in which the best validation accuracy was reached during training. The results of RC and LW are put in Appendix C, available online.

From Table III, we can see that the enhanced method with the robust warm start has significant performance improvements over its original version. The model has better performance when early stopping is not deployed, suggesting that the robust warm start helps not to remember incorrect labels at an early stage. Even after using early stopping, our enhanced version also allows for further performance improvements. Moreover, we presented the results of our ABS method with early stopping...
in the last column. It is usually comparable with the highest accuracy, meaning that our proposal is a simple yet effective baseline for unreliable PLL. The results on CIFAR-10 are also shown in Fig. 1.

VII. CONCLUSION

In this paper, we rethought the forgotten ABS in the era of deep PLL, and improved it theoretically and practically. Theoretically, we proposed two problem settings with different data generation models for noise-free and noisy PLs, and analyzed the conditions that ABS is robust to PLs, which filled the theoretical gap in the robustness analysis of PLL. Practically, we conducted extensive experiments to confirm our theoretical findings, and showed that the IBS methods could be improved from our work, which pushed forward PLL as a whole.

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