Abstract

Label noise has been a practical challenge in deep learning due to the strong capability of deep neural networks in fitting all training data. Prior literature primarily resorts to sample selection methods for combating noisy labels. However, these approaches focus on dividing samples by order sorting or threshold selection, inevitably introducing hyper-parameters (e.g., selection ratio / threshold) that are hard-to-tune and dataset-dependent. To this end, we propose a simple yet effective approach named PNP (Probabilistic Noise Prediction) to explicitly model label noise. Specifically, we simultaneously train two networks, in which one predicts the category label and the other predicts the noise type. By predicting label noise probabilistically, we identify noisy samples and adopt dedicated optimization objectives accordingly. Finally, we establish a joint loss for network update by unifying the classification loss, the auxiliary constraint loss, and the in-distribution consistency loss. Comprehensive experimental results on synthetic and real-world datasets demonstrate the superiority of our proposed method. The source code and models have been made available at https://github.com/NUST-Machine-Intelligence-Laboratory/PNP.

1. Introduction

Although deep neural networks (DNNs) have attained impressive achievements, surpassing traditional methods in various vision tasks [3, 15, 28, 31, 38, 46], their requirement for large-scale high-quality human-labeled training samples (e.g., ImageNet [4] and COCO [21]) can often pose a bottleneck when applied to real-world scenarios. Precise annotation is always labor-expensive and time-consuming, especially when domain-specific expert knowledge is necessary (e.g., fine-grained visual categorization [13, 24, 41]). To alleviate this issue, one promising alternative is to resort to web images for training deep networks [19, 22, 34, 35, 39, 43, 45, 47–49, 53, 54]. However, noisy labels are inevitable in web images [3–4]. It has been demonstrated that noisy labels can impair the performance of deep networks since the over-parameterization equips DNNs with not only large learning capacities but also strong memorization power [16, 52]. Consequently, it is of great significance to develop robust models for learning from noisy labels.

Noisy labels in real-world datasets can be categorized into two types: open-set and closed-set [48]. In the closed-set scenario, the true label of a noisy sample comes from a known label space \( \mathcal{Y}_{\text{known}} \) present in the training data. Conversely, in the open-set scenario, true labels of samples are outside \( \mathcal{Y}_{\text{known}} \). In other words, closed-set noisy samples are in-distribution (ID) ones, while open-set noisy samples are out-of-distribution (OOD) ones. A large body of prior literature primarily focuses on closed-set scenarios,
assumingly, only in-distribution noise exists. However, the ID-noise-only assumption may not hold true in real-world applications. Recently, an increasing number of researchers have been attracted to the open-set noisy problem, which is also the primary focus of this work.

There are mainly two common strategies to tackle noisy labels: loss correction [5, 9, 26, 29, 33, 39, 50, 56] and sample selection [1, 6, 11, 25, 42, 51]. Classic loss correction methods either attempt to estimate the noise transition matrix [2, 5, 9, 26, 33] or seek to regularize losses based on network predictions [29, 56]. Unfortunately, the noise transition matrix is challenging to estimate, while prediction-based loss correction suffers from error accumulation.

Sample selection methods essentially follow an intuitive but straightforward idea: eliminating noisy data and training with the cleaner subset. Researchers have recently witnessed that deep networks tend to fit clean and simple patterns before memorizing noisy labels [16, 52]. Accordingly, many approaches have been proposed to exploit this observation and regard low-loss samples as clean ones. For example, Co-teaching [6] maintains two networks simultaneously and enables them to select low-loss samples for their peer networks. Early sample selection methods usually split samples into two subsets: clean and noisy, neglecting the difference between in-distribution noisy and out-of-distribution noisy labels. More recently, CRSSC [34] and Jo-SRC [48] are proposed to divide samples into three groups: clean ones, in-distribution noisy ones, and out-of-distribution noisy ones, and treat them differently. The former employs a two-step sample selection process to categorize samples into three groups, while the latter proposes global sample selection criteria to distinguish different types of noise. Despite that promising results have been observed, existing methods inevitably involve hard-to-tune and dataset-dependent threshold hyper-parameters for selecting samples, posing a limit to the reliability and scalability of these methods in various larger real-world scenarios.

To address aforementioned issues, we propose a simple yet effective approach, named PNP (Probabilistic Noise Prediction), to probabilistically model label noise in an end-to-end manner. Specifically, we simultaneously train two networks, in which one (i.e., label predictor network) predicts the category of the input data while the other (i.e., noise predictor network) predicts the noise type (i.e., clean / ID noisy / OOD noisy). The clean, ID noisy, OOD noisy samples can be naturally identified according to the prediction from the noise predictor network. To enable effective learning of the noise predictor network, we propose to optimize it in a regression manner, using JS divergence between prediction-label pairs and prediction-prediction pairs. Finally, we impose a consistency regularization on in-distribution data to further advance the learning of our label predictor network and noise predictor network. A comparison between our PNP and existing sample selection methods is visualized in Fig. 1. Our major contributions are:

1. We propose a simple yet effective approach, named PNP, to combat noisy labels. PNP simultaneously predicts the category label and noise type for all training samples. By adopting distinct loss functions for different samples, PNP can robustly learn from noisy training data.

2. PNP employs an auxiliary regression loss for empowering the model to learn to predict the noise type of each sample. JS divergence between prediction-label pairs and prediction-prediction pairs is adopted to approximate the ground-truth noise type. Furthermore, consistency between different views of in-distribution data is encouraged to reinforce the recognition ability.

3. We evaluate two paradigms of sample selection in our method: PNP-hard (hard selection) and PNP-soft (soft selection). We validate the effectiveness and superiority of our method by providing extensive experimental results on both synthetic and real-world noisy datasets. Moreover, comprehensive ablation studies are established to verify each component of our approach.

2. Related Works

Prior works on learning from noisy labels can be briefly categorized into three families:

Label. Early methods primarily focus on correcting corrupted labels. For example, F-correction [26] proposes to adopt a two-step method for estimating the noise transition matrix. S-model [5] proposes to adopt an additional softmax layer to model the noise transition matrix. For these approaches, a well-estimated noise transition matrix is critical in achieving superior and robust performance. However, the noise transition matrix is difficult to estimate, especially in complicated scenarios (e.g., real-world noisy datasets).

Sample. From the perspective of sample, the core idea is to perform sample re-weighting or sample selection. Sample re-weighting methods primarily seek to assign different weights to training samples. For example, Ren et al. [30] propose a meta-learning algorithm to weight training data differently. However, this line of work tends to involve a complicated optimization process and require a small set of clean validation data. Different from sample re-weighting, sample selection methods aim to select correctly-labeled samples for training. Researchers have demonstrated that low-loss samples are more likely to possess correct labels. For example, Co-teaching [6] trains two networks and lets them select low-loss samples for each other. JoCoR [42] employs a joint loss to select low-loss data, encouraging agreement between networks. CRSSC [34] adopts a loss-based selection and a confidence-based selection to identify clean, ID noisy, and OOD noisy samples. Jo-SRC [48] exploits the Jensen-Shannon (JS) divergence and prediction disagreement to globally select different types of noisy data.
Figure 2. The overall framework of PNP. Each input image \( x_i \) is fed into two networks in parallel. The noise predictor network accordingly predicts the probability of \( x_i \) being clean \((P^{\text{clean}})\) / ID \((P^{\text{id}})\) / OOD \((P^{\text{ood}})\). Meanwhile, \( x_i \) is augmented into two weakly and one strongly augmented views before fed into the label predictor network, leading to three label predictions \( p(v_i^c) \), \( p(v_i^w) \), and \( p(v_i^s) \). Afterward, the classification loss \( L_{\text{cls}} \) is computed based on the estimated noise type and the selection paradigm (hard / soft) in the classification module. The constraint loss \( L_{\text{cons}} \) is attained resorting to the approximated ground-truth noise type in the auxiliary module. The consistency loss \( L_{\text{cons}} \) is obtained by encouraging the (label) prediction agreement between different views of in-distribution samples in the consistency module. Finally, our model is updated by back-propagating a joint loss, which is essentially a weighted sum of the above three losses.

3. The Proposed Method

Preliminaries. Given a \( N \)-sample \( C \)-class dataset \( \mathcal{D} = \{(x_i, y_i) | 1 \leq i \leq N\} \), in which \( x_i \) denotes the \( i \)-th training sample and \( y_i \in \{0, 1\}^C \) is its annotated label. We denote the true label of \( x_i \) as \( y_i^t \). Conventionally, we implicitly assume all annotated labels are accurate \( (i.e., y_i = y_i^t) \) and thus optimize the model by minimizing the empirical loss

\[
\mathcal{L} = \mathbb{E}_{\mathcal{D}}[l_{\text{ce}}(x_i, y_i)] = \frac{1}{N} \sum_{i=1}^{N} l_{\text{ce}}(x_i, y_i),
\]

in which

\[
l_{\text{ce}}(x_i, y_i) = - \sum_{c=1}^{C} y_i^c \log(p^c(x_i, \Theta)).
\]

\( \Theta \) denotes the model parameters. \( p^c(x_i, \Theta) \) denotes the predicted softmax probability of the \( i \)-th training sample \( x_i \) over its \( c \)-th class. (For simplicity, we use the notation \( p_i^c \) hereinafter.) Nevertheless, the clean-label assumption may be too restrictive for real-world scenarios and noisy labels are inevitable in many real-world datasets. In this paper, we focus on the scenario where annotated labels are not guaranteed to be correct. Due to the memorization effect [16], noisy labels are prone to leading to inferior performance when used for network training. Thus, it is urgent to design noise-robust methods for addressing noisy labels.

3.1. Probabilistic Noise Modeling

One of the most common strategies for tackling label noise is to find clean samples based on a pre-designed selection process. Owing to the behavior of DNNs in learning simple patterns before fitting noisy labels, previous works have witnessed promising results by selecting low-loss samples as clean ones. However, these methods tend to involve complicated hyper-parameters tuning. For example, Co-teaching [6] and JoCoR [42] require to estimate the noise ratio, while CRSSC [34] and Jo-SRC [48] need to choose a proper selection threshold. Unfortunately, these hyper-parameters \( (e.g., \text{noise ratio and selection threshold}) \) are usually hard to tune and dataset-dependent.

To alleviate the aforementioned issue, we propose to directly model label noise in an end-to-end probabilistic manner. Specifically, we propose to train two parallel networks. The first network, termed as label predictor network (LPN), is trained to predict the category label:

\[
p(x_i) = \sigma(h(f(x_i, \Psi), \Phi_L)) \in \mathbb{R}^C,
\]

where \( \Phi_L \) denotes parameters of the prediction head of LPN. \( \Psi \) denotes parameters of the backbone. \( f(\cdot, \Psi) \) and \( h(\cdot, \Phi_L) \) are mapping functions of the backbone and the prediction head. \( \sigma(\cdot) \) is the softmax function. Conversely, the second network, termed as noise predictor network (NPN), is trained to predict the noise type:

\[
t(x_i) = \sigma(g(f(x_i, \Psi), \Phi_N)) \in \mathbb{R}^3,
\]
in which $\Phi_N$ denotes parameters of the prediction head of NPN. $g(\cdot; \Phi_N)$ is the mapping function of this prediction head. In our implementation, the prediction head of NPN is a multi-layer perception (MLP) network with one hidden layer. Here, we define $t^{(0)}(x_i)$, $t^{(1)}(x_i)$, and $t^{(2)}(x_i)$ as the likelihood of $x_i$ belonging to the clean, ID, and OOD set, respectively. For simplicity, we henceforward denote $P_i^{\text{clean}} = t^{(0)}(x_i)$, $P_i^{\text{id}} = t^{(1)}(x_i)$, and $P_i^{\text{ood}} = t^{(2)}(x_i)$. It should be noted that, in our implementation, to reduce the resource consumption and enable an end-to-end joint optimization, LPN and NPN share the same backbone feature extractor but differ in their prediction heads. By probabilistically modeling label noise, we can conveniently identify and tackle different types of noisy samples accordingly.

### 3.2. Classification Losses for Different Noise

The NPN predicts the noise type of each sample by estimating its “likelihood” of being clean / ID / OOD. We adopt different loss functions for different types of noisy samples. For clean samples, we employ the cross-entropy loss along with an entropy regularization term:

$$l_{\text{clean}}(x_i, y_i) = -\sum_{c=1}^{C} y_i^c \log(p_i^c) - \sum_{c=1}^{C} p_i^c\log(p_i^c).$$

(5)

For in-distribution / out-of-distribution noisy samples, inspired by unsupervised consistency training [43], we propose to treat outputs of strongly and weakly augmented inputs as predictions and targets, respectively. More specifically, for an ID noisy sample $x_i$, we feed its two augmented views (i.e., a strongly augmented one $v_0^i$ and a weakly augmented one $v_w^i$) into our network. The LPN accordingly produces predictions $p(v_0^i)$ and $p(v_w^i)$, which are then leveraged to compute the cross-entropy loss:

$$l_{\text{id}}(x_i) = l_{\text{ce}}(p(v_0^i), \varepsilon(p(v_w^i), \tau)),$$

(6)

in which

$$\varepsilon(z, T) = \frac{\exp(z/T)}{\sum_{z'} \exp(z'/T)}.$$  

(7)

Similarly, for an OOD noisy sample $x_i$, we also employ its two augmented views for computing the classification loss:

$$l_{\text{ood}}(x_i) = l_{\text{ce}}(p(v_0^i), \varepsilon(p(v_w^i), 1/\tau)).$$

(8)

Here, inspired by Jo-SRC [48], we empirically set $\tau = 0.1$, making $\varepsilon(\cdot, \cdot)$ a sharpening operation in Eq. (6) but a flattening operation in Eq. (8).

**Discussion.** The motivation of employing Eqs. (6) and (8) for noisy samples is three-folded. Firstly, by optimizing losses computed from Eqs. (6) and (8), we implicitly enhance consistency between strongly and weakly augmented views of each noisy sample, leading to a smoother model and an improved sample efficiency. Secondly, strong augmentations tend to provide more diverse and natural views, benefiting the generalization performance. Lastly, although first terms (i.e., predictions) in $l_{\text{ce}}$ are identical between Eqs. (6) and (8), the second terms (i.e., targets) are constructed distinctively based on the nature of ID and OOD noisy samples. For ID noisy samples, predictions from a well-trained model tend to be more reliable than given annotations. Therefore, we employ a sharpening operation to advance training by enforcing more confident predictions. On the contrary, OOD noisy samples usually confuse models due to their out-of-task ground-truth categories. By imposing a flattening operation, their predictions will fit an approximately uniform distribution, leading to a boosted robustness and generalization performance.

### 3.3. Constraint of Probabilistic Noise Modeling

We propose to train an additional predictor (i.e., NPN) for estimating the noise type of each sample. However, the NPN is difficult to optimize due to the absence of ground-truth supervision. In this work, we propose to approximate the ground-truth noise type for each sample and accordingly train the NPN. Specifically, we follow Jo-SRC [48] and adopt the Jensen-Shannon (JS) divergence [20] to approximate the probability $Q_i^{\text{clean}}$ of a sample $x_i$ being clean:

$$Q_i^{\text{clean}} = Q_i^{\text{clean}}(x_i) = 1 - D_{\text{JS}}(p(v_0^i)\|y_i).$$

(9)

where $D_{\text{JS}}(\cdot\|\cdot)$ is the JS divergence function. Moreover, inspired by [48], we employ prediction divergence to estimate the “likelihood” $Q_i^{\text{ood}}$ of a sample being OOD. Different from [48], to enable a smoother optimization, we design

$$Q_i^{\text{ood}} = Q_i^{\text{ood}}(x_i) = D_{\text{JS}}(p(v_0^i)\|p(v_w^i)),$$

(10)

in which $v_w^i$ denotes another weakly augmented view of $x_i$.

Once approximations of the ground-truth noise type are obtained, the following auxiliary constraint loss is adopted to optimize the NPN:

$$l_{\text{aux}}(x_i) = |Q_i^{\text{clean}} - Q_i^{\text{clean}}| + |Q_i^{\text{ood}} - Q_i^{\text{ood}}|.$$  

(11)

**Discussion.** (1) Although Eq. (11) only provides a weak constraint due to the ground-truth approximation, the optimization of this auxiliary loss drives the estimation of noise type to its correct direction. (2) The optimization of Eq. (11) is actually a regression task. Therefore, the loss function could be any applicable regression loss (e.g., Mean Absolute Error, Mean Squared Error, etc.). For simplicity, we empirically employ Mean Absolute Error (MAE) loss in our implementation. (3) Jo-SRC [48] uses prediction disagreement to measure the prediction divergence, producing a 0/1 “likelihood”. Conversely, we employ the JS divergence to estimate the prediction disagreement so that our NPN can be optimized in a smoother manner.
3.4. Consistency of In-distribution Data

Intuitively, a well-trained model should predict consistently on different variations of in-distribution samples but contradictorily on those of out-of-distribution data. Due to the employment of prediction divergence in detecting out-of-distribution samples, we propose to impose a consistency regularization loss (i.e., Eq. (12)) on in-distribution data.

\[ l_{cons}(x_i) = D(p(v^w_i)||p(v^w_i)) + D(p(v^w_i)||p(v^w_i)). \] (12)

\( D(\cdot||\cdot) \) denotes the Kullback-Leibler (KL) divergence. The consistency regularization not only implicitly enhances representation learning, but also explicitly empowers our model to better discriminate ID noise and OOD noise.

3.5. PNP-hard and PNP-soft

The overall workflow of our PNP method is shown in Fig. 2. Our algorithm is trained in a two-step manner. Starting with a warm-up period, our network is trained with the original noisy labels by optimizing Eq. (1) for a few epochs. This step facilitates us with a reasonable model for subsequent robust learning. After the warm-up step, we start our PNP training by optimizing the following objective loss function in an end-to-end manner:

\[ \mathcal{L} = \mathcal{L}_{cls} + \gamma \mathcal{L}_{aux} + \omega \mathcal{L}_{cons}, \] (13)

where \( \gamma \) and \( \omega \) are designed to balance different loss terms.

In this work, we evaluate two paradigms of sample selection: hard selection and soft selection. Following the idea of hard sample selection [34,48], PNP-hard employs different loss functions on different types of samples:

\[ \begin{align*}
\mathcal{L}_{cls} &= \mathbb{E}_i [p_i^{clean} \max(p_i^{clean}, p_i^{ood}) I_{clean}(x_i, y_i) + p_i^{ood} I_{ood}(x_i)] \\
\mathcal{L}_{cons} &= \mathbb{E}_i [p_i^{ood} I_{ood}(x_i)].
\end{align*} \] (14)

\( I_A \) is an indicator function, which equals 1 if \( A \) is true, and 0 otherwise. Contrarily, PNP-soft adopts soft sample selection, re-weighting losses based on predictions of noise type:

\[ \begin{align*}
\mathcal{L}_{cls} &= \mathbb{E}_i [p_i^{clean} I_{clean}(x_i, y_i) + p_i^{ood} I_{ood}(x_i)] \\
\mathcal{L}_{cons} &= \mathbb{E}_i [p_i^{ood} I_{ood}(x_i)].
\end{align*} \] (15)

Comparison between PNP-hard and PNP-soft. PNP-hard is intuitive and straightforward, assigning each sample a discrete tag that reveals its noise type. Different loss functions are accordingly employed based on the estimated noise type. On the contrary, PNP-soft adopts a re-weighting schema when computing losses. While hard selection can concretely identify the noise type, it may amplify the risk of incorrect predictions, leading to a potential overfitting problem. Conversely, PNP-soft is beneficial by guaranteeing that at least part of the loss is correctly optimized even if the noise type is wrongly predicted. However, PNP-soft may suffer from the underfitting issue. Empirically, PNP-hard achieves better performance if the noise situation is insignificant and a trustworthy NPN can be attained. When the training data is heavily corrupted, PNP-soft would be superior, owing to its robustness against errors from NPN.

4. Experiments

4.1. Experiment Setup

Datasets. We evaluate our PNP approach on two synthetic datasets (i.e., CIFAR100N and CIFAR80N) and four real-world datasets (i.e., Web-Aircraft, Web-Bird, Web-Car, and Food101N). CIFAR100N and CIFAR80N stem from CIFAR100 [48]. Specifically, we follow Jo-SRC [48] to create the closed-set noisy dataset CIFAR100N and the open-set noisy dataset CIFAR80N. We adopt two classic noise structures: symmetric and asymmetric. Web-Aircraft, Web-Bird, and Web-Car are sub-datasets of WebFG-496 [36], which is a webly supervised fine-grained datasets. Food101N [17] is large-scale real-world noisy dataset.

Evaluation Metric. For assessing the performance of our proposed PNP approach, we adopt the test accuracy as our evaluation metric. Reported results are averaged performance of five repeated experiments under identical settings.

Implementation Details. We adopt a seven-layer DNN [48] for CIFAR100N and CIFAR80N. Adam optimizer [12] is employed during training. We set the initial learning rate as 0.001 and the batch size as 128. We warmup the network for 10 epochs. The learning rate starts to decay linearly after 80 epochs of training. The entire training lasts for 200 epochs. For obtaining further performance gains, we adopt the label smoothing regularization (LSR) [38] technique when calculating clean samples’ classification losses (i.e., Eq. (5)). The LSR parameter \( \epsilon \) is empirically set to 0.6. \( \gamma \) and \( \omega \) are set as 1.0 in default. For Web-Aircraft, Web-Bird, and Web-Car, we leverage ResNet-50 [8] pre-trained on ImageNet as our backbone to compare PNP with other state-of-the-art methods. We update network parameters using SGD optimizer [37] with a momentum of 0.9. The initial learning rate and batch size are 0.0005 and 16, respectively. The warm-up stage lasts for 10 epochs and we train networks for 120 epochs. We start decay learning rate after 10 epochs in a cosine annealing manner. \( \gamma \) and \( \omega \) are also set as 1.0 in default. For Food101N, we follow settings in Jo-SRC [48] and employ pre-trained ResNet-50 for comparison. Default values of \( \gamma \) and \( \omega \) are 1.0 and 0.2.
Baselines. To evaluate our PNP approach on synthetic datasets, we follow Jo-SRC [48] and compare PNP-hard / PNP-soft with state-of-the-art sample selection methods: Decoupling [25], Co-teaching [6], Co-teaching+ [51], Jo-CoR [42], and Jo-SRC [48]. For evaluating on Web-Aircraft, Web-Bird, and Web-Car, we additionally compare PNP with other state-of-the-art methods (e.g., SELFIIE [32], PENCIL [50], AFM [27], CRS2C [34], Self-adaptive [10], DivideMix [18], PLC [55], and Peer-learning [36]). We follow Jo-SRC [48] when evaluating our approach on Food101N. We compare our approach with CleanNet [17], DeepSelf [7], and Jo-SRC [48]. Finally, we denote “Standard” as the baseline case in which we train a deep network using noisy datasets directly. We implement all above methods using PyTorch for performing fair comparison.

4.2. Evaluation on Synthetic Noisy Datasets

We first evaluate PNP on synthetic datasets. By varying the structure and ratio of label noise, we can better understand the effectiveness of PNP in different noise situations.

Results on CIFAR100N. Starting from evaluating our approach in closed-set scenarios, we present the comparison in test accuracy with state-of-the-art approaches on CIFAR100N in Table 1. Results of existing methods are drawn from Jo-SRC [48] and those of our method are obtained under the same experimental settings. From Table 1, we can observe that both PNP-hard and PNP-soft consistently achieve the leading performance. While existing state-of-the-art approaches almost fail in the most inferior case (i.e., Sym-80%), our PNP-hard and PNP-soft still achieve the most appealing performances. We can observe that PNP-hard outperforms PNP-soft only when the noise structure and ratio is Sym-20%. This verifies our argument that hard selection (PNP-hard) will achieve better results only when the noise situation is insignificant. In other cases, PNP-soft consistently performs better than PNP-hard. It should be noted that real-world noisy labels are mostly asymmetric. Table 1 reveals that our PNP-hard / PNP-soft performs notably better than state-of-the-art methods in the case of Asym-40%.

Table 1. Average test accuracy (%) on CIFAR100N and CIFAR80N over the last 10 epochs (“Sym” and “Asym” denote the symmetric and asymmetric label noise, respectively).

<table>
<thead>
<tr>
<th>Methods</th>
<th>CIFAR100N</th>
<th>CIFAR80N</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>Sym – 20%</td>
<td>Sym – 80%</td>
</tr>
<tr>
<td>Standard</td>
<td>35.14 ± 0.44</td>
<td>4.41 ± 0.14</td>
</tr>
<tr>
<td>Decoupling</td>
<td>33.10 ± 0.12</td>
<td>3.89 ± 0.16</td>
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<tr>
<td>Co-teaching</td>
<td>43.73 ± 0.16</td>
<td>15.15 ± 0.46</td>
</tr>
<tr>
<td>Co-teaching+</td>
<td>49.27 ± 0.03</td>
<td>13.44 ± 0.37</td>
</tr>
<tr>
<td>JoCoR</td>
<td>53.01 ± 0.04</td>
<td>15.49 ± 0.98</td>
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<tr>
<td>Jo-SRC</td>
<td>58.15 ± 0.14</td>
<td>23.80 ± 0.05</td>
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<tr>
<td>PNP-hard</td>
<td>64.25 ± 0.12</td>
<td>30.26 ± 0.15</td>
</tr>
<tr>
<td>PNP-soft</td>
<td>63.27 ± 0.14</td>
<td>31.32 ± 0.19</td>
</tr>
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</table>

The remarkable superiority of our method in asymmetric noise indicates that PNP will achieve satisfactory results in real-world noisy datasets.

Results on CIFAR80N. CIFAR80N is specifically created to simulate the real-world (open-set) noisy scenario. The comparison between our method with state-of-the-art approaches is also provided in Tab. 1. Results of existing methods are directly from Jo-SRC [48], and performances of our method are obtained under the same experimental settings. From Tab. 1, we can have the following observations: (1) Our PNP-hard / PNP-soft method consistently outperforms state-of-the-art approaches across different noise scenarios. Our model can achieve the best performance even when facing severe label noise (i.e., Sym-80%). (2) PNP-soft exhibits better performance than PNP-hard in all noisy cases. We believe this results from the complicated noisy labels existed in the open-set noisy dataset CIFAR80N. (3) PNP-hard and PNP-soft obtain impressive performance boost in the case of Asym-40%, validating our design for open-set real-world (asymmetric) problems. These observations firmly validate the effectiveness and superiority of our proposed method in open-set noisy cases.

4.3. Evaluation on Real-world Noisy Datasets

Beyond the above evaluations, we conduct experiments on real-world noisy datasets, including three medium-scale web-image-based fine-grained datasets and one large-scale food dataset, to verify the effectiveness of PNP.

Results on Web-Aircraft / Bird / Car. Web-Aircraft, Web-Bird, and Web-Car are three real-world web image datasets for fine-grained vision categorization. Within each dataset, more than 25% of training samples are associated with unknown (asymmetric) noisy labels. Even worse, these datasets do not provide any label verification information, making it a practical and challenging label noise problem. Tab. 2 illustrates a comparison between our method with state-of-the-art methods. From this table, the leading performance obtained by our method can be witnessed. PNP-hard and PNP-soft both outperform state-of-
Methods Performances (%)

<table>
<thead>
<tr>
<th>Methods</th>
<th>Publications</th>
<th>Backbone</th>
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<th>Web-Bird</th>
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<tr>
<td>AFM</td>
<td>ECCV 2020</td>
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<td>81.04</td>
<td>76.35</td>
<td>83.48</td>
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<td>CRSSC</td>
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<td>81.31</td>
<td>87.68</td>
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<td>Self-adaptive</td>
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<td>78.49</td>
<td>78.19</td>
</tr>
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<td>74.40</td>
<td>84.27</td>
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<tr>
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<td>81.22</td>
<td>88.13</td>
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<td>75.37</td>
<td>82.48</td>
</tr>
<tr>
<td>PNP-hard</td>
<td>-</td>
<td>ResNet50</td>
<td>85.03</td>
<td>81.20</td>
<td>89.93</td>
</tr>
<tr>
<td>PNP-soft</td>
<td>-</td>
<td>ResNet50</td>
<td>85.54</td>
<td>81.93</td>
<td>90.11</td>
</tr>
</tbody>
</table>

Table 2. Comparison with state-of-the-art approaches in test accuracy (%) on Web-Aircraft, Web-Bird, and Web-Car.

Table 3. Comparison with state-of-the-art approaches in test accuracy (%) on Food101N.

<table>
<thead>
<tr>
<th>Method</th>
<th>Backbone</th>
<th>Test accuracy</th>
</tr>
</thead>
<tbody>
<tr>
<td>Standard</td>
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<td>84.51</td>
</tr>
<tr>
<td>CleanNet</td>
<td>ResNet-50</td>
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<td>CleanNet</td>
<td>ResNet-50</td>
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<td>DeepSelf</td>
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<td>Jo-SRC</td>
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<tr>
<td>PNP-hard</td>
<td>ResNet-50</td>
<td>87.31</td>
</tr>
<tr>
<td>PNP-soft</td>
<td>ResNet-50</td>
<td>87.50</td>
</tr>
</tbody>
</table>

Table 4. Impacts of different ingredients in test accuracy (%) on CIFAR80N (Asym-40%). Results at the best epochs are presented.

4.4. Ablation Study

4.4.1 Influence of Different Ingredients

Tab. 4 illustrates impacts of different ingredients in PNP. Clean, ID, and OOD denote the adoption of Eq. (5), Eq. (6), and Eq. (8), respectively. AUX indicates that the constraint loss Eq. (11) is utilized. CONS suggests the employment of in-distribution consistency regularization. LSR is adopted in default. The best result of “Standard” is 29.11%, and that of “Standard + LSR” is 33.10%. From this table, we can observe that each ingredient exhibits a non-trivial significance in our approach. Firstly, by using NPN to identify clean / ID noisy / OOD noisy samples, the performance is promoted by a large margin compared to “Standard + LSR”. Secondly, the employment of the auxiliary constraint empowers the model to achieve more performance boost. Lastly, through imposing consistency regularization on in-distribution data, PNP-hard and PNP-soft are further advanced in robustness.
For studying sensitivity of hyper-parameters, we primarily investigate two parameters (i.e., $\gamma$ and $\omega$) in the value range of $\{0.2, 0.4, 0.6, 0.8, 1.0\}$. Fig. 4 presents the results on the CIFAR80N (Sym-20%) dataset. For better understanding these two hyper-parameters, we additionally provide experimental results of $\gamma = 0$ and $\omega = 0$. The left sub-figure reveals that our method is considerably robust against the value of $\gamma$. The sharp performance increase from $\gamma = 0$ to $\gamma > 0$ demonstrates the importance of employing the auxiliary constraint loss (i.e., Eq. (11)). The right sub-figure exhibits the sensitivity of $\omega$. From this sub-figure, we can observe that while PNP-soft is fairly robust against $\omega$, PNP-hard can benefit from the value increase of this hyper-parameter. Our hypothesis is that the superior noise-robustness of PNP-soft weakens the impact of $\omega$. Since PNP-hard is less robust to label noise, a stronger consistency regularization may better boost the model performance. This sub-figure also reveals a notable performance gap between $\omega = 0$ and $\omega > 0$, manifesting the necessity of adopting in-distribution consistency regularization.

4.4.3 Sensitivity of Hyper-parameters

For studying sensitivity of hyper-parameters, we primarily investigate two parameters (i.e., $\gamma$ and $\omega$) in the value range of $\{0.2, 0.4, 0.6, 0.8, 1.0\}$. Fig. 4 presents the results on the CIFAR80N (Sym-20%) dataset. For better understanding these two hyper-parameters, we additionally provide experimental results of $\gamma = 0$ and $\omega = 0$. The left sub-figure reveals that our method is considerably robust against the value of $\gamma$. The sharp performance increase from $\gamma = 0$ to $\gamma > 0$ demonstrates the importance of employing the auxiliary constraint loss (i.e., Eq. (11)). The right sub-figure exhibits the sensitivity of $\omega$. From this sub-figure, we can observe that while PNP-soft is fairly robust against $\omega$, PNP-hard can benefit from the value increase of this hyper-parameter. Our hypothesis is that the superior noise-robustness of PNP-soft weakens the impact of $\omega$. Since PNP-hard is less robust to label noise, a stronger consistency regularization may better boost the model performance. This sub-figure also reveals a notable performance gap between $\omega = 0$ and $\omega > 0$, manifesting the necessity of adopting in-distribution consistency regularization.

5. Conclusion

In this paper, we focused on the challenge of learning from real-world (open-set) noisy labels. To mitigate their negative impact, we proposed a simple yet effective approach named PNP to model label noise in an end-to-end probabilistic manner. PNP followed the sample selection paradigm but bypassed the requirement for selection thresholds, which were hard-to-tune and dataset-dependent. Specifically, PNP trained two networks in parallel, enabling simultaneous predictions of the category label (i.e., LPN) and the noise type (i.e., NPN). Moreover, a regression task was proposed to optimize the NPN and a consistency regularization was adopted to empower the discrimination ability. Finally, we evaluated two selection paradigms of PNP (i.e., PNP-hard and PNP-soft). A series of experimental results on synthetic and real-world datasets justified the effectiveness and superiority of our proposed approach.

Acknowledgments

This work was supported by National Natural Science Foundation of China (No. 62102182, 61905114, 62072245, and 61932020), Natural Science Foundation of Jiangsu Province (No. BK20210327 and BK20211520), Fundamental Research Funds for the Central Universities (No. 30920021135), and National Key R&D Program of China (No. 2018AAA0102001 and 2021YFF0602101).
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