IF YOUR DATA DISTRIBUTION SHIFTS, USE SELF-LEARNING

ABSTRACT

In this paper, we demonstrate that self-learning techniques like entropy minimization or pseudo-labeling are simple, yet effective techniques for increasing test performance under domain shifts. Our results show that self-learning consistently increases performance under distribution shifts, irrespective of the model architecture, the pre-training technique or the type of distribution shift. At the same time, self-learning is simple to use in practice because it does not require knowledge or access to the original training data or scheme, is robust to hyperparameter choices, is straight-forward to implement and requires only a few training epochs. This makes self-learning techniques highly attractive for any practitioner who applies machine learning algorithms in the real world. We present state-of-the art adaptation results on CIFAR10-C (8.5% error), ImageNet-C (22.0% mCE), ImageNet-R (17.4% error) and ImageNet-A (14.8% error), theoretically study the dynamics of self-supervised adaptation methods and propose a new classification dataset (ImageNet-D) which is challenging even with adaptation.

1 INTRODUCTION

Deep Neural Networks (DNNs) can reach human-level performance in complex cognitive tasks (Brown et al., 2020; He et al., 2016a; Berner et al., 2019) if the distribution of the test data is sufficiently similar to the training data. However, DNNs are known to struggle if the distribution of the test data is shifted relatively to the training data (Geirhos et al., 2018; Dodge & Karam, 2017).

There are two largely distinct communities that aim to increase the performance of models under test-time distribution shifts. The *robustness community* generally considers ImageNet-scale datasets and evaluates models in an *ad-hoc* scenario. Here models are trained on a clean source dataset like ImageNet using, e.g., heavy data augmentation (Hendrycks et al., 2020a; Rusak et al., 2020; Geirhos et al., 2019) and/or large-scale pre-training (Xie et al., 2020; Mahajan et al., 2018). The trained models are not adapted in any way to test-time distribution shifts like corruptions. This evaluation scenario is interesting for applications in which models encounter very different distribution shifts in an unpredictable order (e.g., an online object recognition model), but it misses out on the gains that adaptation to unlabeled samples of the target distribution can bring.

The *Unsupervised Domain Adaptation (UDA) community* often considers smaller-scale datasets and assumes that both the source and the (unlabeled) target dataset are known. Here models are typically trained on both the source and the target dataset (e.g., with an adversarial domain objective, Ganin et al., 2016) before the models are evaluated on a test set with the same distribution as the target. This evaluation scenario provides optimal conditions for adaptation, but the reliance on the source dataset makes methods in this space much more computationally expensive, more impractical and prevents the use of pre-trained models for which the source dataset is unknown or simply too large.

In this work, we consider the *source-free domain adaptation setting*, a middle ground in which models can adapt to the target distribution but without using the source dataset (Kundu et al., 2020; Kim et al., 2021; Li et al., 2020; Liang et al., 2020). This evaluation scenario is interesting for many practitioners and applications. For one, by not being able to use the source dataset, methods in this space need to rely on fine-tuning pre-trained models, which is typically much faster and easier to implement than UDA methods. Second, methods in this space also lend themselves to fast online adaptation in applications where models experience a consistent distribution shift over many inputs.

Figure 1: **A.** Robustness and adaptation to new target datasets has traditionally been achieved by techniques like robust pre-training (with hand-selected or data-driven augmentation strategies), unsupervised domain adaptation (with access to unlabeled samples from the test set), or, more recently, self-supervised learning methods. **B.** We show that on top of these different pre-training tasks, it is always possible (irrespective of architecture, model size or pre-training algorithm) to further adapt models to the target domain with simple self-learning techniques. **C.** We demonstrate these improvements both on small and large-scale datasets, in particular on ImageNet-C, -A, -R and on our new benchmark, ImageNet-D.

Self-learning is part of state-of-the-art UDA methods (French et al., 2018; Shu et al., 2018). We here ask the question whether self-learning methods like *pseudo-labeling* and *entropy-minimization* can also be used in the *source-free* domain adaptation setting. To answer this question, we perform an extensive study of several variants of self-learning, and find consistent and substantial gains in test-time performance across several robustness and out-of-domain benchmarks and a wide range of models and pre-training methods, including models trained with unsupervised domain adapation methods that don't use self-learning. We also find that self-learning outperforms state-of-the-art source-free domain adaptation methods, namely Test-Time Training which is based on a self-supervised auxiliary objective and continual training (Sun et al., 2019b), test-time entropy minimization (Wang et al., 2020) and (gradient-free) BatchNorm adaptation (Schneider et al., 2020; Nado et al., 2020). We perform a large number of ablations to study important design choices for self-learning methods in source-free domain adaptation. Furthermore, we show that a variant of pseudo-labeling with a robust loss function consistently outperforms entropy minimization on ImageNet-scale datasets. We theoretically analyze and empirically verify the influence of the temperature parameter in self-learning and provide guidelines how this single parameter should be chosen. Our approach is visualized in Figure 1.

Our work is conceptually most similar to virtual adversarial domain adaptation performed in the fine-tuning phase of DIRT-T (Shu et al., 2018)) and Test-time entropy minimization (TENT; Wang et al., 2020). In contrast to DIRT-T, we use a simpler adaptation objective, and scale the approach to considerably larger datasets on ImageNet scale. TENT, on the other hand, only evaluated a single method (entropy minimization) on a single vanilla model (ResNet-50) on IN-C. We here substantially expand this analysis to show that self-learning almost universally increases test-time performance under distribution shifts, regardless of the type of distribution shift, the model architecture or the pre-training method.

2 SELF-LEARNING FOR TEST-TIME ADAPTATION

Different variants of self-learning have been used in both unsupervised domain adaptation (French et al., 2018; Shu et al., 2018), self-supervised representation learning (Caron et al., 2021), and in semi-supervised learning (Xie et al., 2020). In a typical self-learning setting a *teacher* network $\mathbf{f}^t$ trained on the source domain predicts labels on the target domain. Then, a second *student* model $\mathbf{f}^s$ is fine-tuned on the predicted labels.

In the following, let $\mathbf{f}^t(\mathbf{x})$ denote the logits for sample $\mathbf{x}$ and let $p^t(j|\mathbf{x}) \equiv \sigma_j(\mathbf{f}^t(\mathbf{x}))$ denote the probability for class j obtained from a softmax function $\sigma_j(\cdot)$. Similarly, $\mathbf{f}^s(\mathbf{x})$ and $p^s(j|\mathbf{x})$ denote the logits and probabilities for the student model $\mathbf{f}^s$. For all techniques, one can optionally

only admit samples where the probability $\max_j p^t(j|\mathbf{x})$ exceeds some threshold. We consider three popular variants of self-learning: Pseudo-labeling with hard or soft labels, as well as entropy minimization.

Hard Pseudo-Labeling (Lee, 2013; Galstyan & Cohen, 2007)   In this setting, one generate labels using a teacher model and train the student model on the pseudo-labels $\tilde{y}$ using the standard cross-entropy loss,

$$\tilde{y} = \mathrm{argmax}_j \, p^t(j|\mathbf{x}), \quad \ell_H(\mathbf{x}, \tilde{y}) := -\log p^s(\tilde{y}|\mathbf{x}). \tag{1}$$

Usually, only samples with a confidence above a certain threshold are considered for training the student. We test several thresholds but note that thresholding means discarding a potentially large portion of the data which leads to a performance decrease in itself. The teacher is updated after each epoch.

Soft Pseudo-Labeling (Lee, 2013; Galstyan & Cohen, 2007)   In contrast to the hard pseudo-labeling variant, we here train the student on class probabilities predicted by the teacher,

$$\ell_S(\mathbf{x}) := -\sum_j p^t(j|\mathbf{x}) \log p^s(j|\mathbf{x}). \tag{2}$$

Soft pseudo-labeling is typically not used in conjunction with thresholding, since it already incorporates the certainty of the model. The teacher is updated after each epoch.

Entropy Minimization (ENT) (Grandvalet & Bengio, 2004)   This variant is similar to soft pseudo-labeling, but we no longer differentiate between a teacher and student network. It corresponds to an "instantaneous" update of the teacher. The training objective becomes

$$\ell_E(\mathbf{x}) := -\sum_j p^s(j|\mathbf{x}) \log p^s(j|\mathbf{x}). \tag{3}$$

Intuitively, self-training with entropy minimization leads to a sharpening of the output distribution for each sample, making the model more confident in its predictions.

Robust Pseudo-Labeling (RPL)   Virtually all introduced self-training variants use the standard cross-entropy classification objective. However, the standard cross-entropy loss has been shown to be sensitive to label noise (Zhang & Sabuncu, 2018; Zhang et al., 2017). In the setting of domain adaptation, inaccuracies in the teacher predictions and, thus, the labels for the student, are inescapable, with severe repercussions for training stability and hyperparameter sensitivity as we show in the results.

As a straight-forward solution to this problem, we propose to replace the cross-entropy loss by a robust classification loss designed to withstand certain amounts of label noise (Ghosh et al., 2017; Song et al., 2020; Shu et al., 2020; Zhang & Sabuncu, 2018). A popular candidate is the *Generalized Cross Entropy (GCE)* loss which combines the noise-tolerant Mean Absolute Error (MAE) loss (Ghosh et al., 2017) with the CE loss. We only consider the hard labels and use the robust GCE loss as the training loss for the student,

$$\ell_{GCE}(\mathbf{x}, \tilde{y}) := q^{-1}(1 - f_{\tilde{y}}^s(\mathbf{x})^q), \tag{4}$$

with $q \in (0, 1]$. For the limit case $q \to 0$, the GCE loss approaches the CE loss and for $q = 1$, the GCE loss is the MAE loss (Zhang & Sabuncu, 2018). We test updating the teacher both after every update step of the student (RPL) and once per epoch (RPL$^{\mathrm{ep}}$).

3   NEW ROBUSTNESS BENCHMARK: IMAGENET-D

Robustness datasets on ImageNet-scale are limited to a few selected domains (image corruptions in IN-C, image renditions in IN-R, difficult images for ResNet50 classifiers in IN-A). In order to test our approach on a wider range of complex distribution shifts, we re-purpose the dataset from the Visual Domain Adaptation Challenge 2019 (DomainNet, Saenko et al., 2019) as an additional robustness benchmark. This dataset comes with six image styles: Clipart, Real, Infograph, Painting, Quickdraw and Sketch. It has 345 classes in total, of which 164 overlap with IN. To benchmark

robustness of IN trained models out of the box, we filter out the classes that cannot be mapped to IN and refer to the smaller version of DomainNet as ImageNet-D (IN-D). We map 463 classes in IN to these 164 IN-D classes, e.g., for an image from the "bird" class in IN-D, we accept all 39 bird classes in IN as valid predictions. We show example images from IN-D in Fig. 2. The detailed evaluation protocol along with justifications for our design choices and additional analysis are outlined in Appendix D.

4 EXPERIMENT DESIGN

Datasets IN-C (Hendrycks & Dietterich, 2019) contains corrupted versions of the 50 000 images in the IN validation set. There are fifteen test and four hold-out corruptions, and there are five severity levels for each corruption. IN-R (Hendrycks et al., 2020a) contains 30 000 images with artistic renditions of 200 classes of the IN dataset. IN-A (an, 2019) is composed of 7500 unmodified real-world images on which standard IN-trained ResNet50 (He et al., 2016b) models yield chance level performance. CIFAR10 (Krizhevsky et al., 2009) and STL10 (Coates et al., 2011) are small-scale image recognition datasets with 10 classes each, and training sets of 50 000/5000 images and test

Figure 2: Example images from the different domains of the new ImageNet-D benchmark.

sets of 10 000/8000 images, respectively. The digit datasets MNIST (Deng, 2012) and MNIST-M (Ganin et al., 2016) both have 60 000 training and 10 000 test images.

Hyperparameters The different self-learning variants have a range of hyperparameters such as the learning rate or the stopping criterion. Our goal is to give a realistic estimation on the performance to be expected in practice.. To this end, we optimize hyperparameters for each variant of pseudo-labeling on a hold-out set of IN-C that contains four types of image corruptions ("speckle noise", "Gaussian blur", "saturate" and "spatter") with five different strengths each, following the procedure suggested in Hendrycks & Dietterich (2019). We refer to the hold-out set of IN-C as our *dev* set.

Models for ImageNet-scale datasets We consider four popular model architectures: ResNet50 (He et al., 2016b), DenseNet161 (Huang et al., 2017), ResNeXt101 (Xie et al., 2017) and EfficientNet-L2 (Tan & Le, 2019) (see Appendix B.1 for details on the used models). For ResNet50, DenseNet and ResNeXt101, we include a simple *vanilla* version trained on IN only. For ResNet50 and ResNeXt101, we additionally include a state-of-the-art robust version trained with DeepAugment and Augmix (DAug+AM, Hendrycks et al., 2020a)[1]. For the ResNeXt model, we also include a version that was trained on 3.5 billion weakly labeled images (IG-3.5B, Mahajan et al., 2018). Finally, for EfficientNet-L2 we select the current state of the art on IN-C which was trained on 300 million images from JFT-300M (Chollet, 2017; Hinton et al., 2014) using a noisy student-teacher protocol (Xie et al., 2020). We validate the IN and IN-C performance of all considered models and match the originally reported scores (Schneider et al., 2020). For EfficientNet-L2, we match IN top-1 accuracy up to 0.1% points, and IN-C up to 0.6% mCE.

Models for CIFAR10/MNIST-scale datasets For CIFAR10-C experiments, we use two WideResNets (WRN, Zagoruyko & Komodakis, 2016): the first one is trained on CIFAR10 and has a depth of 28 and a width of 10 and the second one is trained with AugMix (Hendrycks et al., 2020b) and has a depth of 40 and a width of 2. The remaining small-scale models are trained with unsupervised domain adaptation (UDA) methods. We propose to regard any UDA method which requires joint training with source and target data as a pre-training step, similar to regular pre-training on IN, and use self-learning on top of the final checkpoint. We consider two popular UDA methods: self-supervised domain adaptation (UDA-SS; Sun et al., 2019a) and Domain-Adversarial

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[1]see leaderboard at github.com/hendrycks/robustness

Training of Neural Networks (DANN; Ganin et al., 2016). In UDA-SS, the authors seek to align the representations of both domains by performing an auxiliary self-supervised task on both domains simultaneously. In all UDA-SS experiments, we use a WideResNet with a depth of 26 and a width of 16. In DANN, the authors learn a domain-invariant embedding by optimizing a minimax objective. For all DANN experiments except for MNIST→MNIST-M, we use the same WRN architecture as above. For the MNIST→MNIST-M experiment, the training with the larger model diverged and we used a smaller WideResNet version with a width of 2. We note that DANN training involves optimizing a minimax objective and is generally harder to tune.

All details considering models, and the implementation of the methods can be found in Appendix B.1.

5 RESULTS: SELF-LEARNING UNIVERSALLY IMPROVES MODELS

Self-learning is a powerful learning scheme, and in the following section we show that it allows to perform test-time adaptation on robustified models, models obtained with large-scale pre-training, as well as domain adapted models across a wide range of datasets and distribution shifts. Our main results on large-scale and small-scale datasets are shown in Tables 1 and 2, respectively. These summary tables show final results, and all experiments use the hyperparameters we determined separately on the dev set.

Table 1: Self-learning successfully adapts ImageNet-scale models across different model architectures on IN-C, IN-A and IN-R. We adapt the vanilla ResNet50, ResNeXt101 and DenseNet161 models to IN-C and decrease the mCE by over 19 percent points in all models. Further, self-learning works for models irrespective of their size: Self-learning substantially improves the performance of the ResNet50 and the ResNext101 trained with DAug+AM, on IN-C by 11.9 and 9.7 percent points, respectively. Finally, we further improve the current state-of-the-art model on IN-C—the EfficientNet-L2 Noisy Student model—and report a new state-of-the-art result of 22% mCE (which corresponds to a top1 error of 17.1%) on this benchmark with test-time adaptation (compared to 28% mCE without adaptation).

mCE [%] on IN-C test ($\searrow$)	number of parameters	w/o adapt	w/ adapt RPL	Δ
ResNet50 vanilla (He et al., 2016b)	2.6×10^7	76.7	50.5	(-26.2)
ResNet50 DAug+AM (Hendrycks et al., 2020a)	2.6×10^7	53.6	41.7	(-11.9)
DenseNet161 vanilla (Huang et al., 2017)	2.8×10^7	66.4	47.0	(-19.4)
ResNeXt101$_{32\times8d}$ vanilla (Xie et al., 2017)	8.8×10^7	66.6	43.2	(-23.4)
ResNeXt101$_{32\times8d}$ DAug+AM (Hendrycks et al., 2020a)	8.8×10^7	44.5	34.8	(-9.7)
ResNeXt101$_{32\times8d}$ IG-3.5B (Mahajan et al., 2018)	8.8×10^7	51.7	40.9	(-10.8)
EfficientNet-L2 Noisy Student (Xie et al., 2020)	4.8×10^8	28.3	**22.0**	(-6.3)
top1 error [%] on IN-R ($\searrow$)				
ResNet50 vanilla (He et al., 2016b)	2.6×10^7	63.8	54.1	(-9.7)
EfficientNet-L2 Noisy Student (Xie et al., 2020)	4.8×10^8	23.5	**17.4**	(-6.1)
top1 error [%] on ImageNet-A ($\searrow$)				
EfficientNet-L2 Noisy Student (Xie et al., 2020)	4.8×10^8	16.5	**14.8**	(-1.7)

Self-learning is not limited to the distribution shifts tested in IN-C (like compression artefacts or blur). On IN-R, a dataset with renditions, self-learning improves both the vanilla ResNet50 model and the EfficientNet-L2 model, the latter of which improves from 23.5% to 17.4% top-1 error, a new state of the art. For a vanilla ResNet50, we improve the top-1 error from 63.8% (Hendrycks et al., 2020a) to 54.1%. On IN-A, adapting the EfficientNet-L2 model using self-learning decreases the top-1 error from 16.5% (Xie et al., 2020) to 14.8% top-1 error, again constituting a new state of the art with test-time adaptation on this dataset.

Table 2: Self-learning improves robustified and domain adapted models on small-scale datasets. We test common domain adaptation techniques like DANN (Ganin et al., 2016) and UDA-SS (Sun et al., 2019a), and show that self-learning is effective at further tuning such models to the target domain. We suggest to view unsupervised source/target domain adaptation as a step comparable to pre-training under corruptions, rather than an adaptation technique specifically tuned to the target set—indeed, we can achieve error rates using, e.g., DANN + target adaptation previously only possible with source/target based pseudo-labeling, across different common domain adaptation benchmarks. Self-learning also decreases the error on CIFAR10-C of the Wide ResNet model trained with AugMix (AM, Hendrycks et al., 2020b) and reaches a new state of the art on CIFAR10-C of 8.5% top1 error with test-time adaptation. [†]denotes preliminary results on CIFAR-C dev only, due to instabilities in training the adversarial network in DANN.

top1 error [%] on CIFAR10-C ($\searrow$)	number of parameters	w/o adapt	w/ adapt ENT	Δ
WRN-28-10 vanilla (Zagoruyko & Komodakis, 2016)	3.6×10^7	26.5	12.7	(-13.8)
WRN-40-2 AM (Hendrycks et al., 2020b)	2.2×10^6	11.2	8.5	(-2.7)
WRN-26-16 UDA-SS (Sun et al., 2019a)	9.3×10^7	27.7	16.7	(-11.0)
WRN-26-16 DANN (Ganin et al., 2016)	9.3×10^7	[†]29.7	[†]28.5	(-1.2)
UDA CIFAR10→STL10, top1 error on target [%]($\searrow$)				
WRN-26-16 UDA-SS (Sun et al., 2019a)	9.3×10^7	28.7	22.9	(-5.8)
WRN-26-16 DANN (Ganin et al., 2016)	9.3×10^7	25.0	23.9	(-1.1)
UDA MNIST→MNIST-M, top1 error on target [%]($\searrow$)				
WRN-26-16 UDA-SS (Sun et al., 2019a)	9.3×10^7	4.8	2.0	(-2.8)
WRN-26-2 DANN (Ganin et al., 2016)	1.5×10^6	11.4	5.1	(-6.3)

Table 3: Self-learning also improves large pre-trained models. Unlike BatchNorm adaptation (Schneider et al., 2020), we show that self-learning transfers well to models pre-trained on a large amount of unlabeled data: self-learning decreases the mCE on IN-C of the ResNeXt101 trained on 3.5 billion weakly labeled samples (IG-3.5B, Mahajan et al., 2018) from 51.7% to 40.9%.

mCE on IN-C test [%] ($\searrow$)	no adaptation	BN adaptation	self-learning
ResNeXt101$_{32\times8d}$ vanilla	66.6	56.8	43.2
ResNeXt101$_{32\times8d}$ IG-3.5B	51.7	51.8	**40.9**

Table 4: Self-learning outperforms previously published test-time adaptation approaches on IN-C. The robustness benchmark IN-C has so far mostly been regarded in the ad-hoc evaluation setting as discussed in our introduction. Thus, there are only few published methods that report numbers for test-time adaptation: BatchNorm adaptation (Schneider et al., 2020), Test-Time Training (TTT, Sun et al., 2019b), and TENT (Wang et al., 2020). In particular, note that TTT requires a special loss function at training time, while our approach is agnostic to the pre-training phase. Our self-training results outperforms all three baselines (also after tuning TENT with our full experimental protocol):

mCE on IN-C test [%] ($\searrow$)	w/o adapt	BN adapt	TENT (ours)	self-learning
ResNet50 vanilla	76.7	62.2	53.5 (51.6)	**50.5**
top1 error [%] on IN-C, sev. 5 ($\searrow$)	w/o adapt	BN adapt	TTT	self-learning
ResNet18 vanilla	85.4	72.2	66.3	**61.9**

Table 5: Self-supervised methods based on self-learning allow out-of-the-box test-time adaptation. The recently published DINO method (Caron et al., 2021) is another variant of self-supervised learning that has proven to be effective for unsupervised representation learning. At the core, the method uses soft pseudo-labeling. Here, we test whether a model trained with DINO on the source dataset can be test-time adapted on IN-C using DINO to further improve out-of-distribution performance. Since the used model is a vision transformer model, we test different choices of adaptation parameters and find considerable performance improvements in all cases, yielding an mCE of 43.5%mCE at a parameter count comparable to a ResNet50 model:

mCE on IN-C test [%] ($\searrow$)	w/o adapt	w/ adapt affine layers	w/ adapt bottleneck layers	w/ adapt lin. layers	w/ adapt all weights
ViT-S/16	62.3	51.8	46.8	45.2	**43.5**

6 UNDERSTANDING TEST-TIME ADAPTATION WITH SELF-LEARNING

In the following section, we show ablations and interesting insights of using self-learning for test-time adaptation. If not specified otherwise, all ablations are run on the holdout corruptions of IN-C (our dev set) with a vanilla ResNet50.

Table 6: Robust pseudo-labeling outperforms entropy minimization on large-scale datasets while the reverse is true on small-scale datasets. We find that robust pseudo-labeling consistently improves over entropy minimization on IN-C, while entropy minimization performs better on smaller scale data (CIFAR10, STL10, MNIST). The finding highlights the importance of testing both algorithms on new datasets. The improvement is typically on the order of one percent point:

mCE, IN-C dev	ResNet50	ResNeXt-101	EfficientNet-L2		top-1 err, CIFAR-C	WRN-40
ENT	50.0 ± 0.04	43.0	22.2		ENT	**8.5**
RPL	$\mathbf{48.9} \pm 0.02$	**42.0**	**21.3**		RPL	9.0

Table 7: Robust pseudo-labeling allows usage of the full dataset without a threshold. Classical hard labeling needs a confidence threshold (T) for best performance, thereby reducing the dataset size, while best performance for RPL is reached for full dataset training with a threshold T of 0.0:

diff. self-learning methods	no adapt	soft PL	hard PL (T):	0.0	**0.5**	0.9	RPL (T):	**0.0**	0.5	0.9
mCE on IN-C dev [%]	69.5	60.1		53.8	**51.9**	52.4		**49.7**	49.9	51.8

Table 8: Short update intervals are crucial for fast adaptation. Having established that RPL generally performs better than soft- and hard-labeling, we vary the update interval for the teacher. We find that instant updates are most effective. In entropy minimization, the update interval is instant per default.

Update interval for RPL	w/o adapt	no update	epoch	instant
mCE on IN-C dev [%]	69.5	54.0	49.7	**49.2**

Table 9: Adaptation of only affine layers is important. On IN-C, adapting only the affine parameters works better than adapting all parameters or only the last layer. We indicate the number of adapted parameters in brackets.

Adaptation mechanism	w/o adapt	last layer	full model	affine
mCE on IN-C dev [%]	69.5 [0]	60.2 [2M]	51.5 [22.6M]	**48.9** [5.3K]

Hyperparameters obtained on corruption datasets transfer well to real world datasets. When evaluating models, we select the hyperparameters discussed above (the learning rate and the epoch used for early stopping are the most critical ones) on the holdout set of IN-C. We note that this technique transfers well to IN-R, -A and -D, highlighting the practical value of corruption robustness datasets for adapting models on real distribution shifts.

On IN-D, we performed a control experiment where we selected hyperparameters with leave-one-out cross validation—this selection scheme actually performed worse than IN-C parameter selection (see Appendix D.1).

7 ADAPTING MODELS ON A WIDER RANGE OF DISTRIBUTION SHIFTS REVEALS LIMITATIONS OF ROBUSTIFICATION AND ADAPTATION METHODS

Given the strong performance gains due to self-learning on the tested IN-scale robustness benchmarks IN-C, -R and -A, we are interested how self-learning will perform on a wider range of complex distribution shifts. To test this, we adapt a vanilla ResNet50, several robust IN-C models and the EfficientNet-L2 Noisy Student model on IN-D. We use the same hyperparameters we obtained on IN-C dev for all our IN-D experiments. We show our main results in Table 10.

Table 10: Self-learning decreases the top1 error on some IN-D domains but increases it on others.

domain	Real		Painting		Clipart		Sketch		Infograph		Quickdraw	
adapt	w/o	w/	w/o	w/	w/o	w/	w/o	w/	w/o	w/	w/o	w/
model												
EffNet-L2 Noisy Student	29.2	**27.9**	42.7	**40.9**	45.0	**37.9**	56.4	**51.5**	**77.9**	94.3	**98.4**	99.4
ResNet50 DAug+AM	39.2	36.5	58.7	53.4	68.4	57.0	75.2	61.3	88.1	83.2	98.2	99.1
ResNet50 vanilla	40.1	37.3	65.1	57.8	76.0	63.6	82.0	73.0	89.6	85.1	99.2	99.8

More robust models perform better on IN-D. Comparing the performance of the vanilla ResNet50 model to its robust DAug+AM variant, we find that the DAug+AM model performs better on all domains, with the most significant gains on the "Clipart", "Painting" and "Sketch" domains. We show detailed results for all domains and all tested models in Appendix D.2, along with results on IN-C and IN-R for comparison. We find that the best performing models on IN-D are also the strongest ones on IN-C and IN-R which indicates good generalization capabilities of the techniques combined for these models, given the large differences between the three considered datasets. However, even the best models perform 20 to 30 percentage points worse on IN-D compared to their performance on IN-C or IN-R, indicating that IN-D might be a more challenging benchmark.

All models struggle with some domains of IN-D. The EfficientNet-L2 Noisy Student model obtains the best results on most domains. However, we note that the overall error rates are surprisingly high compared to the model's strong performance on the other considered datasets (IN-A: 14.8% top-1 error, IN-R: 17.4% top-1 error, IN-C: 22.0% mCE). Even on the "Real" domain closest to clean IN where the EfficientNet-L2 model has a top-1 error of 11.6%, the model only reaches a top-1 error of 29.2%. Self-learning decreases the top1 error on all domains except for "Infograph" and "Quickdraw". We note that both domains have very high error rates from the beginning and thus hypothesize that the produced pseudo-labels are of low quality.

Error analysis on IN-D. We investigate the errors a ResNet50 model makes on IN-D by analyzing the most frequently predicted classes for different domains to reveal systematic errors indicative of the encountered distribution shifts. We find most errors interpretable: the classifier assigns the label "comic book" to images from the "Clipart" or "Painting" domains, "website" to images from the "Infograph" domain, and "envelope" to images from the "Sketch" domain. Thus, the classifier predicts the domain rather than the class. We find no systematic errors on the "Real" domain which is expected since this domain should be similar to IN. Detailed results on the top-3 most frequently predicted classes for different domains can be found in Fig. 6, Appendix D.4.

IN-D should be used as an additional robustness benchmark. While the error rates on IN-C, -R and -A are at a well-acceptable level for our largest EfficientNet-L2 model after adaptation, IN-D performance is consistently worse for all models. We propose to move from isolated benchmark settings like IN-R (single domain) to benchmarks more common in domain adaptation (like DomainNet) and make IN-D publicly available as an easy to use dataset for this purpose.

8   A SIMPLE MODEL OF STABILITY IN SELF-LEARNING

We observed that different self-learning schemes are optimal for small-scale vs. large-scale datasets and varying amount of classes. We reconsider the used loss functions, and unify them into

$$\ell(x) = -\sum_j \sigma_j\left(\frac{\mathbf{f}^t(\mathbf{x})}{\tau_t}\right) \log\left(\sigma_j\left(\frac{\mathbf{f}^s(\mathbf{x})}{\tau_s}\right)\right),$$

$$\mathbf{f}^t(\mathbf{x}) = \begin{cases} \mathbf{f}(\mathbf{x}), & \text{entropy minimization} \\ \text{sg}(\mathbf{f}(\mathbf{x})), & \text{pseudo-labeling.} \end{cases}$$

(5)

Here we introduced student and teacher temperature τ_s and τ_t as parameters in the softmax function and the stop gradient operation sg. Caron et al. (2021) fixed τ_s and varied τ_t during

training, and empirically found an upper bound for τ_t above which the training was no longer stable. To better understand such behavior, we study the learning dynamics of the loss function in equation 5 theoretically in a simple two-datapoint, two-classes model with linear student and teacher networks $\mathbf{f}^s(\mathbf{x}) = \mathbf{x}^\top \mathbf{w}^s$ and $\mathbf{f}^t(\mathbf{x}) = \mathbf{x}^\top \mathbf{w}^t$ defined in Appendix A.1. Gradient descent with stop gradient corresponds to hard pseudo-labeling in the limit $\tau_t \to 0$ and to soft pseudo-labeling when $\tau_s = \tau_t = 1$. Gradient descent without stop gradient, i.e., setting $\mathbf{w}^s = \mathbf{w}^t = \mathbf{w}$ corresponds to entropy minimization. We obtain the following result:

Proposition 1 (Collapse in the two-point model)**.** *The student and teacher networks $\mathbf{w}_s$ and $\mathbf{w}_t$ trained with stop gradient collapse to the trivial representation $\forall \mathbf{x}$: $\mathbf{x}^\top \mathbf{w}^s = 0, \mathbf{x}^\top \mathbf{w}^t = 0$ if $\tau_s < \tau_t$. The network $\mathbf{w}$ trained without stop gradient collapses if $\tau_s < \tau_t/2$. Proof. see* § A.2. □

We validate both propositions on a simulated toy model, as well as on the CIFAR-C real dataset and outline the results in Figure 3. In practice, the result suggests that *student temperatures should exceed the teacher temperatures for pseudo-labeling*, and *student temperatures should exceed half the teacher temperature for entropy minimization*. Entropy minimization with standard temperatures ($\tau_s = \tau_t = 1$) and hard pseudo-labeling ($\tau_t \to 0$) are hence stable. The two-point learning dynamics vanish for soft pseudo-labeling with $\tau_s = \tau_t$, suggesting that one would have to analyze a more complex model with more data points. While this does not directly imply that the learning is unstable at this point, we empirically observe that both entropy minimization and hard labeling outperform soft-labeling in practice.

Figure 3: For the two point model, we show accuracy (higher is better) and for the CIFAR10-C simulation, we show the error (lower is better). An important convergence criterion for entropy minimization (top row) and pseudo-labeling (bottom row) is the ratio of student and teacher temperatures; in the two-point model, it lies at $\tau_s = \tau_t$ for PL, and $2\tau_s = \tau_t$ for ENT (first column). Despite the simplicity of the two-point model, the general convergence regions transfer to CIFAR10-C (second column).

9 CONCLUSION

We evaluated and analysed how self-learning, an essential component in many unsupervised domain adaptation and self-supervised pre-training techniques, can be applied for adaptation to both small and large-scale image recognition problems common in robustness research. We demonstrated new state-of-the-art adaptation results with the EfficientNet-L2 model on the benchmarks ImageNet-C, -R, and -A, and introduced a new benchmark dataset (ImageNet-D) which remains challenging even after adaptation. Our theoretical analysis shows the influence of the temperature parameter in the self-learning loss function on the training stability and provides guidelines how to choose a suitable value. We showed that self-learning universally improves test-time performance under diverse distribution shifts irrespective of the architecture or pre-training method. We hope that our work encourages both researchers and practitioners to *use self-learning if their data distribution shifts.*

Reproducility Statement We attempted to make our work as reproducible as possible: We mostly used pre-trained models which are publicly available and we denoted the URL addresses of all used checkpoints; for the checkpoints that were necessary to retrain, we report the Github directories with the source code. We report all used hyperparameters in the Appendix and will release our code upon acceptance of the paper.

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A A TWO-POINT MODEL OF SELF-LEARNING

A.1 DEFINITION OF THE TWO-POINT MODEL

To understand the learning dynamics and properties of different loss functions and their hyperparameters, we propose a simple model of self-learning, both for entropy minimization and pseudo-labeling.

A student network $\mathbf{w}^s \in \mathbb{R}^d$ and a teacher network $\mathbf{w}^t \in \mathbb{R}^d$ are trained on N data points $\{\mathbf{x}_i\}_{i=1}^N$ with the cross-entropy loss function $\mathcal{L}$ defined as

$$\mathcal{L} = -\sum_{i=1}^N \ell(\mathbf{x}_i) = -\sum_{i=1}^N \left(\sigma_t(\mathbf{x}_i^\top \mathbf{w}^t) \log \sigma_s(\mathbf{x}_i^\top \mathbf{w}^s) + \sigma_t(-\mathbf{x}_i^\top \mathbf{w}^t) \log \sigma_s(-\mathbf{x}_i^\top \mathbf{w}^s) \right),$$
(6)

$$\text{where } \sigma_t(z) = \frac{1}{1 + e^{-z/\tau_t}} \text{ and } \sigma_s(z) = \frac{1}{1 + e^{-z/\tau_s}}.$$

Here τ_s and τ_t denote the student and teacher temperature parameters. With stop gradient, student and teacher evolve in time according to

$$\dot{\mathbf{w}}^s = -\nabla_{\mathbf{w}^s} \mathcal{L}\left(\mathbf{w}^s, \mathbf{w}^t\right), \quad \dot{\mathbf{w}}^t = \alpha(\mathbf{w}^s - \mathbf{w}^t),$$
(7)

where α is the learning rate of the teacher. Without stop gradient, student and teacher are set equal to each other, and they evolve as

$$\dot{\mathbf{w}} = -\nabla_{\mathbf{w}} \mathcal{L}(\mathbf{w}), \text{ where } \mathbf{w}^s = \mathbf{w}^t \equiv \mathbf{w}.$$
(8)

We restrict the theoretical analysis to the time evolution of the components of $\mathbf{w}^{s,t}$ in direction of two data points $\mathbf{x}_k$ and $\mathbf{x}_l$, $y_k^{s,t} \equiv \mathbf{x}_k^\top \mathbf{w}^{s,t}$ and $y_l^{s,t} \equiv \mathbf{x}_l^\top \mathbf{w}^{s,t}$. All other components $y_i^{s,t}$ with $i \neq k, l$ are neglected to reduce the dimensionality of the equation system.

$$\text{with stop gradient: } \dot{y}_k^s = -\mathbf{x}_k^\top \nabla_{\mathbf{w}^s} \left(\ell(\mathbf{x}_k) + \ell(\mathbf{x}_l)\right), \quad \dot{y}_l^s = -\mathbf{x}_l^\top \nabla_{\mathbf{w}^s} \left(\ell(\mathbf{x}_k) + \ell(\mathbf{x}_l)\right),$$
$$\dot{y}_k^t = \alpha(y_k^t - y_k^s), \quad \dot{y}_l^t = \alpha(y_l^t - y_l^s),$$
(9)
$$\text{without stop gradient: } \dot{y}_k = -\mathbf{x}_k^\top \nabla_{\mathbf{w}} \left(\ell(\mathbf{x}_k) + \ell(\mathbf{x}_l)\right), \quad \dot{y}_l = -\mathbf{x}_l^\top \nabla_{\mathbf{w}} \left(\ell(\mathbf{x}_k) + \ell(\mathbf{x}_l)\right).$$

A.2 PROOF OF PROPOSITION 1

A.2.1 LEARNING DYNAMICS WITH STOP GRADIENT

Computing the stop gradient evolution defined in equation 7 explicitly yields

$$\dot{\mathbf{w}}^s = -\nabla_{\mathbf{w}^s} \mathcal{L} = \frac{1}{\tau_s} \sum_{i=1}^N \left(\sigma_t(\mathbf{x}_i^\top \mathbf{w}^t)\sigma_s(-\mathbf{x}_i^\top \mathbf{w}^s) - \sigma_t(-\mathbf{x}_i^\top \mathbf{w}^t)\sigma_s(\mathbf{x}_i^\top \mathbf{w}^s) \right) \mathbf{x}_i$$
$$\dot{\mathbf{w}}^t = \alpha(\mathbf{w}^s - \mathbf{w}^t)$$
(10)

The second equality uses the well-known derivative of the sigmoid function, $\partial_z \sigma(z) = \sigma(z)\sigma(-z)$.

The equation system of $2d$ nonlinear, coupled ODEs for $\mathbf{w}^s \in \mathbb{R}^d$ and $\mathbf{w}^t \in \mathbb{R}^d$ in equation 10 is analytically difficult to analyze. Instead of studying the ODEs directly, we act on them with the data points $\mathbf{x}_k^\top$, $k = 1, \ldots, N$, and investigate the dynamics of the components $\mathbf{x}_k^\top \mathbf{w}^{s,t} \equiv y_k^{s,t}$:

$$\dot{y}_k^s = \frac{1}{\tau_s} \sum_{i=1}^N \left(\mathbf{x}_i^\top \mathbf{x}_k\right) \left(\sigma_t(y_i^t)\sigma_s(-y_i^s) - \sigma_t(-y_i^t)\sigma_s(y_i^s) \right)$$
$$\dot{y}_k^t = \alpha(y_k^s - y_k^t).$$
(11)

The learning rate of each mode y_k^s is scaled by $(\mathbf{x}_k^\top \mathbf{x}_i)$ which is much larger for $i = k$ than for $i \neq k$ in high-dimensional spaces. In the two-point approximation, we consider only the two (in absolute value) largest terms $i = k, l$ for a given k in the sum in equation 11. Any changes that $y_k^{s,t}(t)$ and

$y_l^{s,t}(t)$ might induce in other modes $y_i^{s,t}(t)$ are neglected, and so we are left with only four ODEs:

$$
\begin{aligned}
\dot{y}_k^s =& \frac{1}{\tau_s} \|\mathbf{x}_k\|^2 \left(\sigma_t(y_k^t)\sigma_s(-y_k^s) - \sigma_t(-y_k^t)\sigma_s(y_k^s) \right) \\
&+ \frac{1}{\tau_s}(\mathbf{x}_k^\top \mathbf{x}_l) \left(\sigma_t(y_l^t)\sigma_s(-y_l^s) - \sigma_t(-y_l^t)\sigma_s(y_l^s) \right), \\
\dot{y}_l^s =& \frac{1}{\tau_s} \|\mathbf{x}_l\|^2 \left(\sigma_t(y_l^t)\sigma_s(-y_l^s) - \sigma_t(-y_l^t)\sigma_s(y_l^s) \right) \\
&+ \frac{1}{\tau_s}(\mathbf{x}_k^\top \mathbf{x}_l) \left(\sigma_t(y_k^t)\sigma_s(-y_k^s) - \sigma_t(-y_k^t)\sigma_s(y_k^s) \right) \\
\dot{y}_k^t =& \alpha(y_k^s - y_k^t), \ \dot{y}_l^t = \alpha(y_l^s - y_l^t).
\end{aligned}
\tag{12}
$$

The fixed points of equation 12 satisfy

$$
\dot{y}_k^s = \dot{y}_l^s = \dot{y}_k^t = \dot{y}_l^t = 0.
\tag{13}
$$

For $\alpha > 0$, requiring $\dot{y}_k^t = \dot{y}_l^t = 0$ implies that $y_k^s = y_k^t$ and $y_l^s = y_l^t$. For $\tau_s = \tau_t$, the two remaining equations $\dot{y}_k^s = \dot{y}_l^s = 0$ vanish automatically so that there are no non-trivial two-point learning dynamics. For $\tau_s \neq \tau_t$, there is a fixed point at $y_k^{s,t} = y_l^{s,t} = 0$ since at this point, each bracket in equation 12 vanishes individually:

$$
\sigma_t(y_{k,l})\sigma_s(-y_{k,l}) - \sigma_s(-y_{k,l})\sigma_t(y_{k,l}) \Big|_{y_{k,l}=0} = \frac{1}{4} - \frac{1}{4} = 0.
\tag{14}
$$

At the fixed point $y_k^{s,t} = y_l^{s,t} = 0$, $\mathbf{w}^s$ and $\mathbf{w}^t$ are orthogonal to both $\mathbf{x}_k$ and $\mathbf{x}_l$ and so classification fails. If this fixed point is stable, $\mathbf{w}^s$ and $\mathbf{w}^t$ will eventually collapse onto it. The fixed point is stable when all eigenvalues of the Jacobian J of the ODE system equation 12 evaluated at $y_k^{s,t} = y_l^{s,t} = 0$ are negative. This is the case whenever $\tau_s < \tau_t$:

$$
J \Big|_{y_k^{s,t}=y_l^{s,t}=0} = \begin{pmatrix} \frac{\|\mathbf{x}_k\|^2}{4}\left(\frac{1}{\tau_t}-\frac{1}{\tau_s}\right) & \frac{(\mathbf{x}_k^\top \mathbf{x}_l)}{4}\left(\frac{1}{\tau_t}-\frac{1}{\tau_s}\right) & 0 & 0 \\ \frac{(\mathbf{x}_k^\top \mathbf{x}_l)}{4}\left(\frac{1}{\tau_t}-\frac{1}{\tau_s}\right) & \frac{\|\mathbf{x}_l\|^2}{4}\left(\frac{1}{\tau_t}-\frac{1}{\tau_s}\right) & 0 & 0 \\ \alpha & 0 & -\alpha & 0 \\ 0 & \alpha & 0 & -\alpha \end{pmatrix},
$$

eigenvalues: $\lambda_1 = \lambda_2 = -\alpha < 0$,

$$
\lambda_{3,4} = \frac{1}{8}\left(\frac{1}{\tau_t}-\frac{1}{\tau_s}\right)\left(\pm \underbrace{\sqrt{\|\mathbf{x}_k\|^4 + \|\mathbf{x}_l\|^4 - 2\|\mathbf{x}_k\|^2\|\mathbf{x}_l\|^2 + 4(\mathbf{x}_k^\top \mathbf{x}_l)^2}}_{\leq \|\mathbf{x}_k\|^2 + \|\mathbf{x}_l\|^2} + \|\mathbf{x}_k\|^2 + \|\mathbf{x}_l\|^2 \right)
$$

$$
\underbrace{}_{\geq 0 \text{ with equality if } \mathbf{x}_k = \pm \mathbf{x}_l}
\tag{15}
$$

To sum up, training with stop gradient and $\tau_s < \tau_t$ leads to a collapse of the two-point model to the trivial representation $y_k^{s,t} = y_l^{s,t} = 0$.

A.2.2 Learning dynamics without stop gradient

Without stop gradient, we set $\mathbf{w}^t = \mathbf{w}^s \equiv \mathbf{w}$ which leads to an additional term in the gradient:

$$
\begin{aligned}
\dot{\mathbf{w}} = -\nabla_\mathbf{w} \mathcal{L} =& \frac{1}{\tau_s} \sum_{i=1}^N \left(\sigma_t(\mathbf{x}_i^\top \mathbf{w})\sigma_s(-\mathbf{x}_i^\top \mathbf{w}) - \sigma_t(-\mathbf{x}_i^\top \mathbf{w})\sigma_s(\mathbf{x}_i^\top \mathbf{w}) \right)\mathbf{x}_i \\
&+ \frac{1}{\tau_t}\sum_{i=1}^N \sigma_t(\mathbf{x}_i^\top \mathbf{w})\sigma_t(-\mathbf{x}_i^\top \mathbf{w})\underbrace{\left(\log\sigma_s(\mathbf{x}_i^\top \mathbf{w}) - \log\sigma_s(-\mathbf{x}_i^\top \mathbf{w})\right)}_{=\log\left((1+e^{y_i/\tau_s})/(1+e^{-y_i/\tau_s})\right)=y_i/\tau_s}\mathbf{x}_i.
\end{aligned}
\tag{16}
$$

As before in A.2.2, we focus on the evolution of the two components $y_k = \mathbf{w}^\top \mathbf{x}_k$ and $y_l = \mathbf{w}^\top \mathbf{x}_l$.

$$
\begin{aligned}
\dot{y}_k =&\|\mathbf{x}_k\|^2 \left(\frac{1}{\tau_s} \left(\sigma_t(y_k)\sigma_s(-y_k) - \sigma_t(-y_k)\sigma_s(y_k) \right) + \frac{1}{\tau_t}\sigma_t(y_k)\sigma_t(-y_k)y_k \right) \\
&+ (\mathbf{x}_k^\top \mathbf{x}_l) \left(\frac{1}{\tau_s} \left(\sigma_t(y_l)\sigma_s(-y_l) - \sigma_t(-y_l)\sigma_s(y_l) \right) + \frac{1}{\tau_s \tau_t}\sigma_t(y_l)\sigma_t(-y_l)y_l \right) \\
\dot{y}_l =&\|\mathbf{x}_l\|^2 \left(\frac{1}{\tau_s} \left(\sigma_t(y_l)\sigma_s(-y_l) - \sigma_t(-y_l)\sigma_s(y_l) \right) + \frac{1}{\tau_t}\sigma_t(y_l)\sigma_t(-y_l)y_l \right) \\
&+ (\mathbf{x}_k^\top \mathbf{x}_l) \left(\frac{1}{\tau_s} \left(\sigma_t(y_k)\sigma_s(-y_k) - \sigma_t(-y_k)\sigma_s(y_k) \right) + \frac{1}{\tau_s \tau_t}\sigma_t(y_k)\sigma_t(-y_k)y_k \right)
\end{aligned}
\tag{17}
$$

There is a fixed point at $y_k = y_l = 0$ where each bracket in equation 17 vanishes individually,

$$
\frac{1}{\tau_s} \left(\sigma_t(y_{k,l})\sigma_s(-y_{k,l}) - \sigma_t(-y_{k,l})\sigma_s(y_{k,l}) \right) + \frac{1}{\tau_s \tau_t}\sigma_t(y_{k,l})\sigma_t(-y_{k,l})y_{k,l} \Big|_{y_{k,l}} = 0.
\tag{18}
$$

The Jacobian of the ODE system in equation 17 and its eigenvalues evaluated at the fixed point are given by

$$
J\Big|_{y_k=y_l=0} = \begin{pmatrix} \frac{\|\mathbf{x}_k\|^2}{4\tau_s}\left(\frac{2}{\tau_t} - \frac{1}{\tau_s}\right) & \frac{(\mathbf{x}_k^\top \mathbf{x}_l)}{4\tau_s}\left(\frac{2}{\tau_t} - \frac{1}{\tau_s}\right) \\ \frac{(\mathbf{x}_k^\top \mathbf{x}_l)}{4\tau_s}\left(\frac{2}{\tau_t} - \frac{1}{\tau_s}\right) & \frac{\|\mathbf{x}_l\|^2}{4\tau_s}\left(\frac{2}{\tau_t} - \frac{1}{\tau_s}\right) \end{pmatrix}
$$

$$
\lambda_{1,2} = \frac{1}{8\tau_s}\left(\frac{2}{\tau_t} - \frac{1}{\tau_s}\right)\underbrace{\left(\pm \underbrace{\sqrt{\|\mathbf{x}_k\|^4 + \|\mathbf{x}_l\|^4 - 2\|\mathbf{x}_k\|^2\|\mathbf{x}_l\|^2 + 4(\mathbf{x}_k^\top \mathbf{x}_l)^2}}_{\leq \|\mathbf{x}_k\|^2 + \|\mathbf{x}_l\|^2} + \|\mathbf{x}_k\|^2 + \|\mathbf{x}_l\|^2 \right)}_{\geq 0 \text{ with equality if } \mathbf{x}_k = \pm \mathbf{x}_l}.
\tag{19}
$$

Hence the fixed point is stable and the two-point model without stop gradient collapses to $y_k = y_l = 0$ whenever $\tau_s < \tau_t/2$.

B   ADDITIONAL INFORMATION ON USED MODELS

B.1   DETAILS ON ALL HYPERPARAMETERS WE TESTED FOR DIFFERENT MODELS

For all models except EfficientNet-L2, we adapt the batch norm statistics to the test domains following (Schneider et al., 2020). We do not expect significant gains for combining EfficientNet-L2 with batch norm adaptation: as demonstrated in (Schneider et al., 2020), models trained with large amounts of weakly labeled data do not seem to benefit from batch norm adaptation.

ResNet50 models   We use a vanilla ResNet50 model and compare soft- and hard-labeling against entropy minimization and robust pseudo-labeling. To find optimal hyperparameters for all methods, we perform an extensive evaluation and test (i) three different adaptation mechanisms (ii) several learning rates 1.0×10^{-4}, 1.0×10^{-3}, 1.0×10^{-2} and 5.0×10^{-2}, (iii) the number of training epochs and (iv) updating the teacher after each epoch or each iteration. For all experiments, we use a batch size of 128. The hyperparameter search is performed on IN-C dev. We then use the optimal hyperparameters to evaluate the methods on the IN-C test set.

ResNeXt101 models   The ResNeXt101 model is considerably larger than the ResNet50 model and we therefore limit the number of ablation studies we perform for this architecture. Besides a baseline, we include a state-of-the-art robust version trained with DeepAugment+Augmix (DAug+AM, Hendrycks et al., 2020a) and a version that was trained on 3.5 billion weakly labeled images (IG-3.5B, Mahajan et al., 2018). We only test the two leading methods on the ResNeXt101 models (ENT and RPL). We vary the learning rate in same interval as for the ResNet50 model but scale it down linearly to account for the smaller batch size of 32. We only train the affine batch normalization parameters because adapting only these parameters leads to the best results

on ResNet50 and is much more resource efficient than adapting all model parameters. Again, the hyperparameter search is performed only on the development corruptions of IN-C. We then use the optimal hyperparameters to evaluate the methods on the IN-C test set.

EfficientNet-L2 models   The current state of the art on IN, IN-C, IN-R and IN-A is an EfficientNet-L2 trained on 300 million images from JFT-300M (Chollet, 2017; Hinton et al., 2014) using a noisy student-teacher protocol (Xie et al., 2020). We adapt this model for only one epoch due to resource constraints. During the hyperparameter search, we only evaluate three corruptions on the IN-C development set[2] and test the learning rates 4.6×10^{-2}, 4.6×10^{-3}, 4.6×10^{-4} and 4.6×10^{-5}. We use the optimal hyperparameters to evaluate ENT and RPL on the full IN-C test set (with all severity levels).

UDA-SS models   We trained the models using the scripts from the official code base at github.com/yueatsprograms/uda_release. We used the provided scripts for the cases: (a) source: CIFAR10, target: STL10 and (b) source: MNIST, target: MNIST-M. For the case (c) source: CIFAR10, target: CIFAR10-C, we used the hyperparameters from case (a) since this case seemed to be the closest match to the new setting. We think that the baseline performance of the UDA-SS models can be further improved with hyperparameter tuning.

DANN models   To train models with the DANN-method, we used the PyTorch implementation of this paper at https://github.com/fungtion/DANN_py3. The code base only provides scripts and hyperparameters for the case (b) source: MNIST, target: MNIST-M. For the cases (a) and (c), we used the same optimizer and trained the model for 100 epochs. We think that the baseline performance of the DANN models can be further improved with hyperparameter tuning.

Preprocessing   For IN, IN-R, IN-A and IN-D, we resize all images to 256×256 px and take the center 224×224 px crop. The IN-C images are already rescaled and cropped. We center and re-scale the color values with $\mu_{RGB} = [0.485, 0.456, 0.406]$ and $\sigma_{RGB} = [0.229, 0.224, 0.225]$. For the EfficientNet-L2, we follow the procedure in Xie et al. (2020) and rescale all inputs to a resolution of 507×507 px and then center-crop them to 475×475 px.

B.2   FULL LIST OF USED MODELS

ImageNet scale models   ImageNet trained models (ResNet50, DenseNet161, ResNeXt) are taken directly from torchvision (Marcel & Rodriguez, 2010). The model variants trained with DeepAugment and AugMix augmentations (Hendrycks et al., 2020b;a) are taken from https://github.com/hendrycks/imagenet-r. The weakly-supervised ResNeXt101 model is taken from the PyTorch Hub. For EfficientNet (Tan & Le, 2019), we use the PyTorch re-implementation available at https://github.com/rwightman/gen-efficientnet-pytorch. This is a verified re-implementation of the original work by Xie et al. (2020). We verify the performance on ImageNet, yielding a 88.23% top-1 accuracy and 98.546% top-5 accuracy which is within 0.2% points of the originally reported result (Xie et al., 2020). On ImageNet-C, our reproduced baseline achieves 28.9% mCE vs. 28.3% mCE originally reported by Xie et al. (2020). As noted in the re-implementation, this offset is possible due to minor differences in the pre-processing. It is possible that our adaptation results would improve further when applied on the original codebase by Xie *et al.*.

Small scale models   We train the UDA-SS models using the original code base at github.com/yueatsprograms/uda_release, with the hyperparameters given in the provided bash scripts. For our DANN experiments, we use the PyTorch implementation at github.com/fungtion/DANN_py3. We use the hyperparameters in the provided bash scripts.

The following Table contains all models we evaluated on various datasets with references and links to the corresponding source code.

---

[2]We compare the results of computing the dev set on the 1, 3 and 5 severities versus the 1, 2, 3, 4 and 5 severities on our ResNeXt101 model in the Supplementary material.

Model	Source
WideResNet(28,10) (Croce et al., 2020)	https://github.com/RobustBench/robustbench/tree/master/robustbench
WideResNet(40,2)+AugMix (Croce et al., 2020)	https://github.com/RobustBench/robustbench/tree/master/robustbench
ResNet50 (He et al., 2016b)	https://github.com/pytorch/vision/tree/master/torchvision/models
ResNeXt101, 32×8d (He et al., 2016b)	https://github.com/pytorch/vision/tree/master/torchvision/models
DenseNet (Huang et al., 2017)	https://github.com/pytorch/vision/tree/master/torchvision/models
ResNeXt101, 32×8d (Xie et al., 2017)	https://pytorch.org/hub/facebookresearch_WSL-Images_resnext/
ResNet50+DeepAugment+AugMix (Hendrycks et al., 2020a)	https://github.com/hendrycks/imagenet-r
ResNext101 (Hendrycks et al., 2020a)	https://github.com/hendrycks/imagenet-r
ResNext101 32×8d IG-3.5B (Mahajan et al., 2018)	https://github.com/facebookresearch/WSL-Images/blob/master/hubconf.p
Noisy Student EfficientNet-L2 (Xie et al., 2020)	https://github.com/rwightman/gen-efficientnet-pytorch

Table 11: mCE in % on the IN-C dev set for ENT and RPL for different numbers of training epochs when adapting the affine batch norm parameters of a ResNet50 model.

Table 12: mCE ($\searrow$) in % on the IN-C dev set for different learning rates for EfficientNet-L2. We favor $q = 0.8$ over $q = 0.7$ due to slightly improved robustness to changes in the learning rate in the worst case error setting.

criterion		ENT			RPL	
lr	1e-4	1e-3	1e-2	1e-4	1e-3	1e-2
epoch						
0	60.2	60.2	60.2	60.2	60.2	60.2
1	54.3	**50.0**	72.5	57.4	51.1	52.5
2	52.4	50.9	96.5	55.8	49.6	57.4
3	51.5	51.0	112.9	54.6	49.2	64.2
4	51.0	52.4	124.1	53.7	49.0	71.0
5	50.7	53.5	131.2	52.9	**48.9**	76.3
6	50.7	53.5	131.2	52.9	48.9	76.3

lr (4.6 ×)	base	10^{-3}	10^{-4}	10^{-5}	10^{-6}
ENT	25.5	87.8	25.3	**22.2**	24.1
RPL$_{q=0.7}$	25.5	60.3	**21.3**	23.3	n/a
RPL$_{q=0.8}$	25.5	58.2	**21.4**	23.4	n/a

Table 13: mCE in % on IN-C dev for entropy minimization for different learning rates and training epochs for ResNeXt101. (div.=diverged)

Table 14: mCE in % on IN-C dev for robust pseudo-labeling for different learning rates and training epochs for ResNeXt101. (div.=diverged)

ENT	Baseline			IG-3.5B			DAug+AM		
lr 2.5 ×	1e-4	1e-3	5e-3	1e-4	1e-3	5e-3	1e-4	1e-3	5e-3
epoch									
BASE	53.6	53.6	53.6	47.4	47.4	47.4	37.4	37.4	37.4
1	**43.0**	92.2	div.	40.9	40.4	58.6	**35.4**	46.4	div.
2	44.8	118.4	div.	39.8	41.5	69.5	35.5	90.8	div.
3	45.4	131.9	div.	39.3	42.6	76.1	35.5	122.5	div.
4	46.7	div.	div.	**39.1**	44.2	84.3	35.6	133.8	div.

RPL	Baseline			IG-3.5B			DAug+AM		
lr 2.5×	1e-4	1e-3	5e-3	1e-4	1e-3	5e-3	1e-4	1e-3	5e-3
epoch									
BASE	53.6	53.6	53.6	47.4	47.4	47.4	37.4	37.4	37.4
1	43.4	51.3	div.	45.0	39.9	43.6	35.3	35.1	79.1
2	42.3	63.2	div.	43.4	**39.3**	48.2	34.9	35.6	121.2
3	42.0	72.6	div.	42.4	39.4	52.9	34.7	40.1	133.5
4	**42.0**	72.6	div.	42.4	39.4	52.9	**34.7**	40.1	133.5

C   DETAILED AND ADDITIONAL RESULTS ON IN-C

C.1   DETAILED RESULTS FOR TUNING EPOCHS AND LEARNING RATES

We tune the learning rate for all models and the number of training epochs for all models except the EfficientNet-L2. In this section, we present detailed results for tuning these hyperparameters for all considered models. The best hyperparameters that we found in this analysis, are summarized in Table 15.

Table 15: The best hyperparameters for all models that we found on IN-C. For all models, we fine-tune only the affine batch normalization parameters and use $q = 0.8$ for RPL. The small batchsize for the EfficientNet model is due to hardware limitations.

Model	Method	Learning rate	batch size	number of epochs
vanilla ResNet50	ENT	1×10^{-3}	128	1
vanilla ResNet50	RPL	1×10^{-3}	128	5
vanilla ResNeXt101	ENT	2.5×10^{-4}	128	1
vanilla ResNeXt101	RPL	2.5×10^{-4}	128	4
IG-3.5B ResNeXt101	ENT	2.5×10^{-4}	128	4
IG-3.5B ResNeXt101	RPL	2.5×10^{-3}	128	2
DAug+AM ResNeXt101	ENT	2.5×10^{-4}	128	1
DAug+AM ResNeXt101	RPL	2.5×10^{-4}	128	4
EfficientNet-L2	ENT	4.6×10^{-5}	8	1
EfficientNet-L2	RPL	4.6×10^{-4}	8	1

Figure 4: Severity-wise mean corruption error (normalized using the *average* AlexNet baseline error for each corruption) for ResNet50 (RN50), ResNext101 (RNx101) variants and the Noisy Student L2 model. Especially for more robust models (DeepAugment+Augmix and Noisy Student L2), most gains are obtained across higher severities 4 and 5. For weaker models, the baseline variant (Base) is additionally substantially improved for smaller corruptions.

C.2 DETAILED RESULTS FOR ALL IN-C CORRUPTIONS

We outline detailed results for all corruptions and models in Table 16. Performance across the severities in the dataset is depicted in Figure 4. All detailed results presented here are obtained by following the model selection protocol outlined in the main text.

C.3 ABLATION OVER THE HYPERPARAMETER q FOR RPL

For RPL, we must choose the hyperparameter q. We performed an ablation study over q and show results in Table 17, demonstrating that RPL is robust to the choice of q, with slight preference to higher values. Note: In the initial parameter sweep for this paper, we only compared $q = 0.7$ and $q = 0.8$. Given the result in Table 17, it could be interesting to re-run the models in Table 1 of the main paper with $q = 0.9$, which could yield another (small) improvement in mCE.

q	0.5	0.6	0.7	0.8	0.9
mCE (dev)	49.5	49.3	49.2	49.2	49.1

Table 17: ImageNet-C dev set mCE in %, vanilla ResNet50, batch size 96. We report the best score across a maximum of six adaptation epochs.

C.4 SELF-TRAINING OUTPERFORMS CONTRASTIVE TEST-TIME TRAINING (SUN ET AL., 2019B)

Sun et al. (2019b) use a ResNet18 for their experiments on ImageNet and only evaluate their method on severity 5 of IN-C. To enable a fair comparison, we trained a ResNet18 with both hard labeling and RPL and compare the efficacy of both methods to Test-Time Training in Table 18. For both hard labeling and RPL, we use the hyperparameters we found for the vanilla ResNet50 model and thus, we expect even better results for hyperparameters tuned on the vanilla ResNet18 model and following our general hyperparameter search protocol.

While all methods (self-learning and TTT) improve the performance over a simple vanilla ResNet18, we note that even the very simple baseline using hard labeling already outperfoms Test-Time Training; further gains are possible with RPL. The result highlights the importance of simple baselines (like self-learning) when proposing new domain adaptation schemes. It is likely that many established DA techniques more complex than the basic self-learning techniques considered in this work will even further improve over TTT and other adaptation approaches developed exclusively in robustness settings.

C.5 EFFECT OF BATCH SIZE AND LINEAR LEARNING RATE SCALING

How is self-learning performance affected by batch size constraints? We compare the effect of different batch sizes and linear learning rate scaling. In general, we found that affine adaptation

Table 16: Detailed results for each corruption along with mean corruption error (mCE) as reported in Table 2 in the main paper. We show (unnormalized) top-1 error rate averaged across 15 test corruptions along with the mean corruption error (mCE: which is normalized). Hyperparameter selection for both ENT and RPL was carried out on the dev corruptions as outlined in the main text. Mismatch in baseline mCE for EfficientNet-L2 can be most likely attributed to pre-processing differences between the original tensorflow implementation Xie et al. (2020) and the PyTorch reimplementation we employ. We start with slightly weaker baselines for ResNet50 and ResNext101 than Schneider et al. (2020): ResNet50 and ResNext101 results are slightly worse than previously reported results (typically 0.1% points) due to the smaller batch size of 128 and 32. Smaller batch sizes impact the quality of re-estimated batch norm statistics when computation is performed on the fly Schneider et al. (2020), which is of no concern here due to the large gains obtained by pseudo-labeling.

	gauss	shot	impulse	defocus	glass	motion	zoom	snow	frost	fog	bright	contrast	elastic	pixelate	jpeg	**mCE**
ResNet50																
Baseline Schneider et al. (2020)																62.2
Baseline (ours)	57.2	59.5	60.0	61.4	62.3	51.3	49.5	54.6	54.1	39.3	29.1	46.7	41.4	38.2	41.8	62.8
ENT	45.5	45.5	46.8	48.4	48.7	40.0	40.3	42.0	46.6	33.2	28.1	42.4	35.2	32.2	35.1	51.6
RPL	44.2	44.4	45.5	47.0	47.4	38.8	39.2	40.7	46.2	32.5	27.7	42.7	34.6	31.6	34.4	50.5
ResNeXt101 Baseline																
Baseline Schneider et al. (2020)																56.7
Baseline (ours)	52.8	54.1	54.0	55.4	56.8	46.7	46.6	48.5	49.4	36.6	25.4	42.8	37.8	32.5	36.7	56.8
ENT	40.5	39.5	41.4	41.6	43.0	34.1	34.5	35.0	39.4	28.5	24.0	33.8	30.3	27.2	30.5	44.3
RPL	39.4	38.9	39.8	40.3	41.0	33.4	33.8	34.6	38.7	28.0	23.7	31.4	29.8	26.8	30.0	43.2
ResNeXt101 IG-3.5B																
Baseline Schneider et al. (2020)																51.6
Baseline (ours)	50.7	51.5	53.1	54.2	55.5	45.5	44.7	41.7	42.0	28.1	20.1	33.8	35.4	27.8	33.9	51.8
ENT	38.6	38.3	40.4	41.4	41.5	33.8	33.6	32.2	34.6	24.1	19.7	26.3	27.6	24.2	27.9	40.8
RPL	39.1	39.2	40.8	42.1	42.4	33.7	33.5	31.8	34.7	23.9	19.6	26.1	27.5	23.8	27.5	40.9
ResNeXt101 DeepAug+Augmix																
Baseline Schneider et al. (2020)																38.0
Baseline (ours)	30.0	30.0	30.2	32.9	35.5	28.9	31.9	33.3	32.8	29.5	22.6	28.4	31.2	23.0	26.5	38.1
ENT	28.7	28.5	29.0	29.8	30.9	26.9	28.0	29.3	30.5	26.2	23.2	26.3	28.5	23.7	26.0	35.5
RPL	28.1	27.8	28.3	29.1	30.1	26.3	27.4	28.8	29.8	25.9	22.7	25.6	27.9	23.2	25.4	34.8
Noisy Student L2																
Baseline Xie et al. (2020)																28.3
Baseline (ours)	21.6	22.0	20.5	23.9	40.5	19.8	23.2	22.8	26.9	21.0	15.2	21.2	24.8	17.9	18.6	28.9
ENT	18.5	18.7	17.4	18.8	23.4	16.9	18.8	17.1	19.6	16.8	14.1	16.6	19.6	15.8	16.5	23.0
RPL	17.8	18.0	17.0	18.1	21.4	16.4	17.9	16.4	18.7	15.7	13.6	15.6	19.2	15.0	15.6	22.0

Table 18: Comparison to Test-Time Training Sun et al. (2019b): Top-1 error for a ResNet18 and severity 5 for all corruptions.

	gauss	shot	impulse	defocus	glass	motion	zoom	snow	frost	fog	bright	contrast	elastic	pixelate	jpeg	**Avg**
vanilla ResNet18	98.8	98.2	99.0	88.6	91.3	88.8	82.4	89.1	83.5	85.7	48.7	96.6	83.2	76.9	70.4	85.41
Test-Time Training	73.7	71.4	73.1	76.3	93.4	71.3	66.6	64.4	81.3	52.4	41.7	**64.7**	55.7	52.2	55.7	66.26
hard labeling, single epoch	73.2	70.8	73.6	76.5	75.6	63.9	56.1	59.0	**65.9**	48.4	39.7	85.2	50.4	47.0	51.5	62.45
RPL (ours), four epochs	**71.3**	**68.3**	**71.7**	76.2	75.6	**61.5**	**54.4**	**56.9**	67.1	**47.3**	**39.3**	93.2	**48.9**	**45.7**	50.4	**61.85**

experiments on ResNet50 scale can be run with batch size 128 on a Nvidia V100 GPU (16GB), while only batch size 96 experiments are possible on RTX 2080 GPUs.

The results in Table 19 show that for a ResNet50 model, higher batch size yields a generally better performance.

Table 19: ImageNet-C dev set mCE for various batch sizes with linear learning rate scaling. All results are computed for a vanilla ResNet50 model using RPL with $q = 0.8$, reporting the best score across a maximium of six adaptation epochs.

batch size	16	32	64	80	96	128
learning rate ($\times 10^{-3}$)	0.125	0.250	0.500	0.625	0.750	1
dev mCE	53.8	51.0	49.7	49.3	49.2	48.9

Table 20: mDE in % on IN-D for different model selection strategies.

model	model selection	
	L1outCV	IN-C dev
ResNet50 RPL$_{q=0.8}$	81.3	76.1
ResNet50 ENT	82.4	77.3
EfficientNet-L2 ENT	69.2	66.8
EfficientNet-L2 RPL$_{q=0.8}$	69.1	67.2

Table 21: Top-1 error on IN-D in % as obtained by robust ResNet50 models. For reference, we also show the mCE on IN-C and the top-1 error on IN-R.

Model	Clipart	Infograph	Painting	Quickdraw	Real	Sketch	mDE	IN-C	IN-R
vanilla	76.0	89.6	65.1	99.2	40.1	82.0	88.2	76.7	63.9
SIN Geirhos et al. (2019)	71.3	88.6	62.6	97.5	40.6	77.0	85.6	69.3	58.5
ANT Rusak et al. (2020)	73.4	88.9	63.3	99.2	39.9	80.8	86.9	62.4	61.0
ANT+SIN Rusak et al. (2020)	68.4	88.6	60.6	95.5	40.8	70.3	83.1	60.7	53.7
AugMix Hendrycks et al. (2020b)	70.8	88.6	62.1	99.1	39.0	78.5	85.4	65.3	58.9
DeepAugment Hendrycks et al. (2020a)	72.0	88.8	61.4	98.9	39.4	78.5	85.6	60.4	57.8
DeepAug+Augmix Hendrycks et al. (2020a)	68.4	88.1	58.7	98.2	39.2	75.2	83.4	53.6	53.2

D DETAILED AND ADDITIONAL RESULTS ON IN-D

D.1 EVALUATION PROTOCOL ON IN-D

The domains in IN-D differ in terms of their difficulty for the studied models. Therefore, to calculate an aggregate score, we propose normalizing the error rates by the error achieved by AlexNet on the respective domains to calculate the mean error, following the approach in Hendrycks & Dietterich (2019) for IN-C. This way, we obtain the aggregate score mean Domain Error (mDE) by calculating the mean over different domains,

$$\text{DE}_d^f = \frac{E_d^f}{E_d^{\text{AlexNet}}}, \qquad \text{mDE} = \frac{1}{D}\sum_{d=1}^{D} E_d^f, \qquad (20)$$

where E_d^f is the top-1 error of a classifier f on domain d.

Leave-one-out-cross-validation For all IN-D results we report in this paper, we chose the hyperparameters on the IN-C dev set. We tried a different model selection scheme on IN-D as a control experiment with "Leave one out cross-validation" (L1outCV): with a round-robin procedure, we choose the hyperparameters for the test domain on all other domains. We select the same hyperparameters as when tuning on the "dev" set: For the ResNet50 model, we select over the number of training epochs (with a maximum of 7 training epochs) and search for the optimal learning rate in the set [0.01, 0.001, 0.0001]. For the EfficientNet-L2 model, we train only for one epoch as before and select the optimal learning rate in the set [4.6×10^{-3}, 4.6×10^{-4}, 4.6×10^{-5}, 4.6×10^{-6}]. This model selection leads to worse results both for the ResNet50 and the EfficientNet-L2 models, highlighting the robustness of our model selection process, see Table 20.

D.2 DETAILED RESULTS FOR ROBUST RESNET50 MODELS ON IN-D

We show detailed results for all models on IN-D for vanilla evaluation (Table 21) BN adaptation (Table 22), RPL$_{q=0.8}$ (Table 23) and ENT(Table 24). For RPL$_{q=0.8}$ and ENT, we use the same hyperparameters that we chose on our IN-C 'dev' set. This means we train the models for 5 epochs with RPL$_{q=0.8}$ and for one epoch with ENT.

The summary results for all models are shown in Table 25.

We show the top-1 error for the different IN-D domains versus training epochs for a vanilla ResNet50 in Fig. 5. We indicate the epochs 1 and 5 at which we extract the errors with dashed black lines.

Table 22: Top1 error on IN-D in % as obtained by state-of-the-art robust ResNet50 models and batch norm adaptation, with a batch size of 128.

Model	Clipart	Infograph	Painting	Quickdraw	Real	Sketch	mDE
vanilla	70.2	88.2	63.5	97.8	41.1	78.3	80.2
SIN Geirhos et al. (2019)	67.3	89.7	62.2	97.2	44.0	75.2	79.6
ANT Rusak et al. (2020)	69.2	89.4	63.0	97.5	42.9	79.5	80.7
ANT+SIN Rusak et al. (2020)	64.9	88.2	60.0	96.8	42.6	73.0	77.8
AugMix Hendrycks et al. (2020b)	66.9	88.1	61.2	97.1	40.4	75.0	78.4
DeepAugment Hendrycks et al. (2020a)	66.6	89.7	60.0	97.2	42.5	75.1	78.8
DeepAug+Augmix Hendrycks et al. (2020a)	61.9	85.7	57.5	95.3	40.2	69.2	74.9

Table 23: Top-1 error on IN-D in % as obtained by state-of-the-art robust ResNet50 models and $\text{RPL}_{q=0.8}$.

Model	Clipart	Infograph	Painting	Quickdraw	Real	Sketch	mDE
vanilla	63.6	85.1	57.8	99.8	37.3	73.0	76.1
SIN Geirhos et al. (2019)	60.8	86.4	56.0	99.0	37.8	67.0	76.8
ANT Rusak et al. (2020)	63.4	86.3	57.7	99.2	37.7	71.0	78.1
ANT+SIN Rusak et al. (2020)	61.5	86.4	56.8	97.0	39.0	67.1	76.1
AugMix Hendrycks et al. (2020b)	59.7	83.4	54.1	98.2	35.6	70.1	74.6
DeepAugment Hendrycks et al. (2020a)	58.1	84.6	53.3	99.0	36.2	64.2	74.8
DeepAug+Augmix Hendrycks et al. (2020a)	57.0	83.2	53.4	99.1	36.5	61.3	72.6

Figure 5: Top-1 error for the different IN-D domains for a ResNet50 and training with $\text{RPL}_{q=0.8}$ and ENT. We indicate the epochs at which we extract the test errors by the dashed black lines (epoch 1 for ENTand epoch 5 for $\text{RPL}_{q=0.8}$).

Table 24: Top-1 error on IN-D in % as obtained by state-of-the-art robust ResNet50 models and ENT.

Model	Clipart	Infograph	Painting	Quickdraw	Real	Sketch	mDE
vanilla	65.1	85.8	59.2	98.5	38.4	75.8	77.3
SIN Geirhos et al. (2019)	62.1	87.0	57.3	99.1	39.0	68.6	75.5
ANT Rusak et al. (2020)	64.2	86.9	58.7	97.1	38.8	72.8	76.5
ANT+SIN Rusak et al. (2020)	62.2	86.8	57.7	95.8	40.1	68.7	75.2
AugMix Hendrycks et al. (2020b)	60.2	84.6	55.8	97.6	36.8	72.0	74.4
DeepAugment Hendrycks et al. (2020a)	59.5	85.7	54.4	98.0	37.1	66.4	73.3
DeepAug+Augmix Hendrycks et al. (2020a)	58.4	84.3	54.7	98.5	38.1	63.6	72.7

Table 25: mDE on IN-D in % as obtained by robust ResNet50 models with a baseline evaluation, batch norm adaptation, $\text{RPL}_{q=0.8}$ and ENT.

| Model | mDE on IN-D ($\searrow$) | | | |
	Baseline	BN adapt	$\text{RPL}_{q=0.8}$	ENT
vanilla	88.2	80.2	76.1	77.3
SIN Geirhos et al. (2019)	85.6	79.6	76.8	75.5
ANT Rusak et al. (2020)	86.9	80.7	78.1	76.5
ANT+SIN Rusak et al. (2020)	**83.1**	77.8	76.1	75.2
AugMix Hendrycks et al. (2020b)	85.4	78.4	74.6	74.4
DAug Hendrycks et al. (2020a)	85.6	78.8	74.8	73.3
DAug+AM Hendrycks et al. (2020a)	83.4	**74.9**	72.6	**72.7**

Table 26: Top-1 error ($\searrow$) on IN-D in % for EfficientNet-L2

Domain	Baseline	ENT	RPL
Clipart	45.0	39.8	**37.9**
Infograph	**77.9**	91.3	94.3
Painting	42.7	41.7	**40.9**
Quickdraw	**98.4**	99.4	99.4
Real	29.2	28.7	**27.9**
Sketch	56.4	**48.0**	51.5
mDE	67.2	**66.8**	67.2

D.3 DETAILED RESULTS FOR THE EFFICIENTNET-L2 NOISY STUDENT MODEL ON IN-D

We show the detailed results for the EfficientNet-L2 Noisy Student model on IN-D in Table 26.

D.4 DETAILED RESULTS ON THE ERROR ANALYSIS ON IN-D

Frequently predicted classes We analyze the most frequently predicted classes on IN-D by a vanilla ResNet50 and show the results in Fig. 6. We make several interesting observations: First, we find most errors interpretable: it makes sense that a ResNet50 assigns the label "comic book" to images from the "clipart" or "painting" domains, or "website" to images from the "infograph" domain, or "envelope" to images from the "sketch" domain. Second, on the hard domain "quickdraw", the ResNet50 mostly predicts non-sensical classes that are not in IN-D, mirroring its almost chance performance on this domain. Third, we find no systematic errors on the "real" domain which is expected since this domain should be similar to IN.

Filtering predictions on IN-D that cannot be mapped to ImageNet-D We perform a second analysis: We filter the predicted labels according to whether they can be mapped to IN-D and report the filtered top-1 errors as well as the percentage of filtered out inputs in Table 27. We note that for the domains "infograph" and "quickdraw", the ResNet50 predicts labels that cannot be mapped to IN-D in over 70% of all cases, highlighting the hardness of these two domains.

Filtering labels and predictions on IN that cannot be mapped to ImageNet-D To test for possible class-bias effects, we test the performance of a ResNet50 model on IN classes that can be mapped to IN-D and report the results in Table 27.

First, we map IN labels to IN-D to make the setting as similar as possible to our experiments on IN-D and report the top-1 error (12.1%). This error is significantly lower compared to the top-1 error a ResNet50 obtains following the standard evaluation protocol (23.9%). This can be explained by the simplification of the task: While in IN there are 39 bird classes, these are all mapped to the same hierarchical class in IN-D. Therefore, the classes in IN-D are more dissimilar from each other than in IN. Additionally, there are only 164 IN-D classes compared to the 1000 IN classes, raising the chance level prediction.

If we further only accept predictions that can be mapped to IN-D, the top-1 error is slightly increased to 13.4%. In total, about 52.7% of all images in the IN validation set cannot be mapped to IN-D.

Figure 6: Systematic predictions of a vanilla ResNet50 on IN-D for different domains.

Table 27: top-1 error on IN and different IN-D domains for different settings: left column: default evaluation, middle column: predicted labels that cannot be mapped to IN-D are filtered out, right column: percentage of filtered out labels.

Dataset	top-1 error in %	top-1 error on filtered labels in %	percentage of rejected inputs
IN val	12.1	13.4	52.7
IN-D real	40.2	17.2	27.6
IN-D clipart	76.1	59.0	59.0
IN-D infograph	89.7	59.3	74.6
IN-D painting	65.2	39.5	42.4
IN-D quickdraw	99.3	96.7	76.1
IN-D sketch	82.1	65.6	47.9

D.5 TOP-1 ERROR ON IN-D FOR ALEXNET

We report the top-1 error numbers on different IN-D as achieved by AlexNet in Table 28. We used these numbers for normalization when calculating mDE.

Table 28: top-1 error on IN-D by AlexNet which was used for normalization.

Dataset	top-1 error in %
IN-D real	54.887
IN-D clipart	84.010
IN-D infograph	95.072
IN-D painting	79.080
IN-D quickdraw	99.745
IN-D sketch	91.189

E SOFTWARE STACK

We use different open source software packages for our experiments, most notably Docker (Merkel, 2014), scipy and numpy (Virtanen et al., 2020), GNU parallel (Tange, 2011), Tensorflow (Abadi et al., 2016), PyTorch (Paszke et al., 2017), and torchvision (Marcel & Rodriguez, 2010).