Training-Free Uncertainty Estimation for Dense Regression: Sensitivity as a Surrogate

Lu Mi 1  Hao Wang 2  Yonglong Tian 1  Hao He 1  Nir Shavit 1

Abstract

Uncertainty estimation is an essential step in the evaluation of the robustness for deep learning models in computer vision, especially when applied in risk-sensitive areas. However, most state-of-the-art deep learning models either fail to obtain uncertainty estimation or need significant modification (e.g., formulating a proper Bayesian treatment) to obtain it. Most previous methods are not able to take an arbitrary model off the shelf and generate uncertainty estimation without retraining or redesigning it. To address this gap, we perform a systematic exploration into training-free uncertainty estimation for dense regression, an unrecognized yet important problem, and provide a theoretical construction justifying such estimations. We propose three simple and scalable methods to analyze the variance of outputs from a trained network under tolerable perturbations: infer-transformation, infer-noise, and infer-dropout. They operate solely during the inference, without the need to re-train, re-design, or fine-tune the models, as typically required by state-of-the-art uncertainty estimation methods. Surprisingly, even without involving such perturbations in training, our methods produce comparable or even better uncertainty estimation when compared to training-required state-of-the-art methods.

1. Introduction

Deep neural networks have achieved remarkable or even super-human performance in many tasks (Krizhevsky et al., 2012; He et al., 2015; Silver et al., 2016). While most previous work in the field has focused on improving accuracy in various tasks, in several risk-sensitive areas such as autonomous driving (Chen et al., 2015) and healthcare (Zhang et al., 2019), reliability and robustness are arguably more important and interesting than accuracy.

Recently, several novel approaches have been proposed to take into account an estimation of uncertainty during training and inference (Huang et al., 2018). Some use probabilistic formulations for neural networks (Graves, 2011; Hernández-Lobato & Adams, 2015; Wang et al., 2016; Shekhovtsov & Flach, 2018) and model the distribution over the parameters (weights) and/or the neurons. Such formulations naturally produce distributions over the possible outputs (Ilg et al., 2018; Yang et al., 2019). Others utilize the randomness induced during training and inference (e.g., dropout and ensembling) to obtain an uncertainty estimation (Gal & Ghahramani, 2016; Lakshminarayanan et al., 2017; Kendall et al., 2015).

All methods above require specific designs or a special training pipeline in order to involve the uncertainty estimation during training. Unfortunately, there are many cases where such premeditated designs or pipelines cannot be implemented. For example, if one wants to study the uncertainty of trained models released online, retraining is not always an option, especially when only a black-box model is provided or the training data is not available. Moreover, most models are deterministic and do not have stochasticity. A straightforward solution is to add dropout layers into proper locations and finetune the model (Gal & Ghahramani, 2016). However, this is impractical for many state-of-the-art and published models, especially those trained on large datasets (e.g. ImageNet (Deng et al., 2009)) with a vast amount of industrial computing resources. In addition, models that have already been distilled, pruned, or binarized fall short of fitting re-training (Han et al., 2015; Hou et al., 2016).

To fill this gap, we identify the problem of training-free uncertainty estimation: how to obtain an uncertainty estimation of any given model without re-designing, re-training, or fine-tuning it. We focus on two scenarios: black-box uncertainty estimation (BBUE), where one has access to the model only as a black box, and gray-box uncertainty estimation (GBUE), where one has access to intermediate-layer neurons of the model (but not the parameters). Our work is a systematic exploration into this unrecognized yet
important problem.

In our paper, we focus on regression tasks. We note that for classification tasks, the softmax output is naturally a distribution. Methods that use entropy for uncertainty estimation qualify as a training-free method and have outperformed MC-Dropout (Bahat & Shakhnarovich, 2018; Gal & Ghahramani, 2016; Hendrycks & Gimpel, 2016; Wang et al., 2019) (see Appendix for experiment results). Regression tasks are more challenging than classification problems since there is no output distribution. Our major contributions are:

1. We perform a systematic exploration of training-free uncertainty estimation for regression models and provide a theoretical construction justifying such estimations.
2. We propose simple and scalable methods, infer-transformation, infer-noise, and infer-dropout, using a tolerable perturbation to effectively and efficiently estimate uncertainty.
3. Surprisingly, we find that our methods are able to generate uncertainty estimation comparable or even better than training-required baselines in real-world large-scale dense regression tasks.

2. Methodology

Three Cases on Parameter Accessibility. We distinguish among three cases based on accessibility of the original model. 1. Black-box case: the model is given as a trained black box without any access to its internal structure. 2. Gray-box case: the internal representations (feature maps) of the model is accessible (while the parameters are not) and can be modified during inference. 3. White-box case: the model is available for all modifications (e.g., its weights can be modified, which requires training). In this paper we focus on the black-box and gray-box cases, for which we offer, correspondingly, two classes of methods in Fig. 1.

2.1. Black-Box Uncertainty Estimation: Infer-Transformation

Given a black-box model, we explore the behavior of the outputs for different transformed versions of the input. Specifically, we transform the input with tolerable perturbations, e.g., perturbations that do not cause significant increase in the loss (see Sec. 2.3 for details), and then use the variance of the perturbed outputs as estimated uncertainty. Here we focus on transformations that preserve pertinent characteristics of the input, such as rotations, flips, etc. Formally, given an input image X, our measured uncertainty is defined as $\mathbb{V}[Z] = \mathbb{V}_{T}[T' \circ F \circ T(X)]$, where $T \in T$ is a transformation, $T'$ is $T$’s inverse operation, and $F$ is a function representing the black-box neural network. $Z = T' \circ F \circ T(X)$ is a sample from the perturbed output distribution. Note that it is possible to sample $Z = F(X)$, where $P$ happens to be a 360-degree rotation.

2.2. Gray-Box Uncertainty Estimation: Infer-Noise and Infer-Dropout

Given a gray-box model, we consider another class of methods for generating multiple outputs from a distribution: randomly perturbing latent codes. Compared with the black-box case, this provides finer granularity on modulating the perturbation strength to ensure tolerability. Specifically we propose infer-noise, which introduces Gaussian noise at an intermediate layer of the trained model, and infer-dropout, which uses dropout instead. For infer-noise, the noise will be added to the feature maps of a certain layer. This noise is randomly sampled multiple times during inference to form a set of diverse outputs. For infer-dropout, random dropout is performed for multiple forwards to generate output samples, the variance of which are then used as uncertainty estimation. Formally, given an input image X, our measured uncertainty is defined as $\mathbb{V}[Z] = \mathbb{V}_{P}[F_2 \circ P \circ F_1(X)]$, where $P$ is sampled from a perturbation set $P$ (e.g., Gaussian noise with $\sigma = 1$). $F_1$ is the function of network layers before the perturbation $P$, $F_2$ represents network layers after $P$, and $F_2 \circ F_1(X)$ is the gray-box network $F(X)$. $Z = F_2 \circ P \circ F_1(X)$ is a sample from the perturbed output distribution. Note that it is possible to sample $Z = F(X)$, where $P$ happens to be a perturbation noise of all zeros.

2.3. Sensitivity as a Surrogate Measure

The idea at the core of our approach is to impose tolerable perturbations on the original trained model’s input or
intermediate representations (feature maps). Given a perturbation output sample \( Z \) and ground truth \( Y \), the perturbation tolerability is defined using \( C = |E[Z] - Y| \leq \epsilon \), where \( \epsilon \) denotes the perturbation error threshold; smaller \( \epsilon \) indicates better tolerability. Such perturbations generate a sensitivity map \( V[Z] \) as a surrogate measure of the model’s uncertainty. Note that our method involves no sacrifice in the predictive performance of original model. One can use our method to produce uncertainty estimation while still use the original model to make predictions.

**Lemma 2.1** (Chebyshev’s Inequality). Let \( Z \) be any random variable with variance \( V[Z] < \infty \). Then for a constant margin \( t \geq 0 \), \( \mathbb{P}(|Z - E[Z]| \geq t) \leq V[Z]/t^2 \).

**Variance and Sensitivity.** If \( Z \) is the model prediction, the probability \( \mathbb{P}(|Z - E[Z]| \geq t) \) translates to ‘how possible the model prediction \( Z \) deviates from \( E[Z] \) by a margin larger than \( t \)’ and therefore measures the ‘model sensitivity’. Such sensitivity is bounded by a scaled variance \( V[Z]/t^2 \) (where \( t^2 \) is a constant). Therefore, Lemma 2.1 connects our output variance \( V[Z] \) to model sensitivity.

**Theorem 2.1.** Let \( Y \) be the ground truth and \( Z \) be a random variable representing the model prediction, where randomness comes from our perturbation. We have that

\[
\mathbb{P}(|Z - Y| \geq t) \leq V[Z]/(t - C)^2,
\]

for any constant margin \( t \geq C \), where \( C = |E[Z] - Y| \) is the prediction error for \( E[Z] \).

**Variance and Uncertainty.** Similar to Lemma 2.1, \( \mathbb{P}(|Z - Y| \geq t) \) in Theorem 2.1 translates to ‘how possible the model prediction \( Z \) deviates from the ground truth \( Y \) by a margin larger than \( t \)’ and therefore measures the ‘uncertainty’. Theorem 2.1 establishes \( V[Z]/(t - C)^2 \) as an upper bound for the ‘uncertainty’. In practice, we directly use \( V[Z] \) since \( C \) is the unknown ground-truth prediction error. In Sec. 3, we empirically show that \( V[Z] \) as a rough approximation for \( V[Z]/(t - C)^2 \) can already obtain uncertainty estimation on par with or even better than state-of-the-art baselines. Moreover, to eliminate the effect of \( t \), one could integrate \( t \) out on both sides and get the summary of ‘uncertainty’ over different thresholds \( t' \).

\[
\int_{t=0}^{\infty} \mathbb{P}(|Z - Y| \geq t) dt \leq C + 2\sqrt{V[Z]}.
\]

We can see that the variance is an upper bound for such a summary of ‘uncertainty’ over different thresholds.

**Why We Need Tolerability.** (1) Wider valid region: The upper bound in Eqn. 1 is valid only when \( t \geq C \); therefore since \( C \leq \epsilon \) by definition, better tolerability (i.e. lower \( \epsilon \)) leads to smaller \( C \), giving the bound a wider valid region w.r.t. the margin \( t \). (2) Tighter bound: Better tolerability also guarantees smaller \( C \) and \( V[Z] \), making \( V[Z]/(t - C)^2 \) smaller and consequently a tighter bound given a constant margin \( t \) (see Sec. 3.1 and Fig. 3 for details). From Eqn. 2, we can also see that smaller \( C \) leads to a tighter bound on the summary uncertainty.

**Epistemic Uncertainty.** The uncertainty estimated in our approach is epistemic uncertainty. Assuming the model perfectly fits infinite data – all variants of augmented “data” (including both data inputs and intermediate features) applied with different perturbations – we will get zero variance, meaning zero epistemic uncertainty.

### 3. Experiments

In this section, we evaluate our three proposed approaches in two representative real-world large-scale dense regression tasks, super resolution and depth estimation. More results
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Table 1. Mean/patch-wise/block-wise/pixel-wise AUSE and correlation between $L_1$ loss and uncertainty, and NLL on SR benchmark dataset Set 14. Our infer-transformation, infer-dropout and infer-noise are compared with MC-dropout (Gal & Ghahramani, 2016), deep ensemble (Lakshminarayanan et al., 2017) and log likelihood maximization (LLM) (Zhang et al., 2019). MC-drop$^1$ uses the output of the original model as prediction while MC-drop$^2$ uses the mean of output samples from the re-trained model (with added dropout) as prediction. Models evaluated: SRGAN.

Figure 3. Visualization of probability $P(|Z - Y| > t)$ (solid) and the upper bound $\mathbb{V}[Z]/(t - C)^2$ (dashed) versus $t$ for three representative pixels, empirically verifying the validity of $\mathbb{V}[Z]/(t - C)^2$ as the uncertainty’s upper bound. Note that as the perturbation strength (i.e. dropout rate) gets larger, both $\mathbb{V}[Z]$ and $C$ increase and consequently loosen the bound (as mentioned in Sec. 2.3); reflected in Fig. 3, we can see the area between the probability curve and the bound curve of each pixel also gets larger from Fig. 3 (left) to Fig. 3 (right), which leads to worse performance (lower average pixel-wise correlation). This highlights the need for tolerable perturbations (i.e. low $\epsilon$ and $C$ in Sec. 2.3). Note that this is also consistent with Eqn. 2 since the area under each curve (in the valid region $t > C$) corresponds to the integration over $t$. Table 4 in Appendix shows correlation and AUSE for methods compared to the oracle (performance bound) $\mathbb{V}[Z]/(t - C)^2$. Our methods, especially infer-transformation, are reasonably close to the oracle and compare favorably with baselines.

The Role of Tolerable Perturbations. Tolerable perturbations play a crucial role in obtaining effective uncertainty estimation. Better tolerability means smaller decrease in accuracy after perturbation (i.e. smaller $\epsilon$ in Sec. 2.3). Fig. 5 in Appendix shows the optimal cases to generate uncertainty maps with high correlation require that $C$ should remain small after perturbation (high tolerability).

3.1. Experiment Results

Qualitative Results. Fig. 2 shows some qualitative results for an example image in the SR task. We can see that the variance maps generated in our task are consistent to the level of ambiguity. Specifically, in our methods, high variance occurs in areas with high randomness and high frequency. As expected, these high-variance areas usually correspond to large prediction error.

Comparative Performance with Training-required Baselines. We report correlation, AUSE and NLL using the optimal hyper-parameters in different methods in Table 1. Meanwhile, we report the evaluation on other tasks in Table 2 and Table 3 in Appendix. Based on these metrics, our methods infer-transformation, infer-dropout and infer-noise could provide comparable or even better results than the training-required state-of-the-art baselines. Even a small number of samples are able to guarantee sufficient quality. For the super-resolution task, we find infer-transformation achieves the highest performance and even outperforms training-required baselines. For depth estimation, infer-noise outperforms other baselines.

related to depth estimation are in Appendix.
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A. Proof of Theorem 3.1.

Proof. Replacing $t$ in Lemma 3.1 with $(t - C)$ where $t \geq C$, we have $P(|Z - \mathbb{E}[Z]| + C \geq t) \leq \mathbb{V}[Z]/(t - C)^2$, meaning that $P(|Z - \mathbb{E}[Z]| + C < t) > \mathbb{V}[Z]/(t - C)^2$. Notice the triangle inequality:

$$|Z - Y| \leq |Z - \mathbb{E}[Z]| + |\mathbb{E}[Z] - Y| = |Z - \mathbb{E}[Z]| + C.$$  

We therefore have $P(|Z - Y| < t) > \mathbb{V}[Z]/(t - C)^2$, which is equivalent to $P(|Z - Y| \geq t) \leq \mathbb{V}[Z]/(t - C)^2$, completing the proof. □

B. Related Work

Probabilistic Neural Networks for Uncertainty Estimation. Probabilistic neural networks consider the input and model parameters as random variables which take effect as the source of stochasticity (Nix & Weigend, 1994; Welling & Teh, 2011; Graves, 2011; Hernández-Lobato & Adams, 2015; Wang et al., 2016). Traditional Bayesian neural networks model the distribution over the parameters (weights) (MacKay, 1992; Hinton & Van Camp, 1993; Graves, 2011; Welling & Teh, 2011) and obtain the output distribution by marginalizing out the parameters. Even with recent improvement (Balan et al., 2015; Hernández-Lobato & Adams, 2015), one major limitation is that the size of network at least doubles under this assumption, and the propagation with a distribution is usually computationally expensive. Another set of popular and efficient methods (Gal & Ghahramani, 2016; Teye et al., 2018) formulate dropout (Srivastava et al., 2014) or batch normalization (Ioffe & Szegedy, 2015) as approximations to Bayesian neural networks. For example, MC-dropout (Gal & Ghahramani, 2016) injects dropout into some layers during both training and inference (Tsymbalov et al., 2019). Unlike most models that disable dropout during inference, MC-dropout feed-forwards the same example multiple times with dropout enabled, in order to form a distribution on the output. Meanwhile, other works (Wang et al., 2016; Shekhovtsov & Flach, 2018) propose sampling-free probabilistic neural networks as a lightweight Bayesian treatment for neural networks.

Non-probabilistic Neural Networks for Uncertainty Estimation. Other strategies (Zhao et al., 2020) such as deep ensemble (Lakshminarayanan et al., 2017; Huang et al., 2017a; Ashukha et al., 2020) train an ensemble of neural networks from scratch, where some randomness is induced during the training process, i.e. the initial weight is randomly sampled from a distribution. During inference, these networks will generate a distribution of the output. Though simple and effective, training multiple networks costs even more time and memory than Bayesian neural networks. Another efficient method log likelihood maximization (LLM) is to train the network to have both original outputs and uncertainty predictions, by jointly optimizing both (Zhang et al., 2019; Poggi et al., 2020). Besides the methods above focusing on uncertainty in classification models; there are also works investigating uncertainty in regression models (Kuleshov et al., 2018; Song et al., 2019; Zelikman et al., 2020). However, all methods above requires re-training, introduces heavy implementation overhead, and sometimes makes the optimization process more challenging.

C. Experiment details

In this section, we evaluate our three proposed approaches in two representative real-world large-scale dense regression tasks, super resolution and depth estimation.

C.1. Single Image Super Resolution

The task of Single Image Super Resolution (SR) is to reconstruct a high-resolution (HR) image from a low-resolution (LR) input. Here we focus on analyzing the state-of-the-art SRGAN model (Ledig et al., 2017), which can restore photo-realistic high-quality images. SRGAN always outputs deterministic restorations since the conditional GAN (Mirza & Osindero, 2014) used in this model involves no latent variable sampling. However, we can still evaluate its uncertainty with our proposed methods.

We apply our methods to estimate uncertainty in one open-source version of this work (Dong et al., 2017). The package provides two models trained with different loss functions: 1) SRResnet model with $L_2$ loss and 2) SRGAN model with a combination of $L_2$ loss and adversarial loss. We evaluate our methods on both models in the black-box(gray-box settings.

Infer-Transformation. For infer-transformation, we apply rotation of $K \times 90$ degrees ($K = 0, 1, 2, 3$) as well as horizontal flip to the LR input, feed it into the trained model during the inference, and apply the inverse transformation to its output. We could generate at most 8 samples using this strategy, and then calculate the pixel-wise variance.

Infer-Noise. In infer-noise, we take the trained model and add a Gaussian-noise layer, which has standard deviation $\sigma \in \{0.01, 0.02, 0.05, 0.1, 0.2, 0.5\}$ and mean 0, at different locations (layers). We choose 4 different locations for noise injection, including the layers right after the input and some intermediate layers (see details in Sec. F). For each experiment, we only add the noise into one layer with a specific $\sigma$ value. Sample numbers of 8 and 32 are evaluated.

Infer-Dropout. In infer-dropout, we take the trained model and add a dropout layer with varied dropout rates. We choose the dropout rate $\rho$ from the set $\{0.01, 0.02, 0.05, 0.1, 0.2, 0.5\}$ and use the same set of locations as the infer-noise. For each experiment, we only add the layer into one location with one specific dropout rate.
We evaluate the model on NYU Depth Dataset V2. For inference, we train these networks with the same number of epochs until they converge. During inference, each of them generates a single deterministic output, with 4 or 8 samples generated in total. The third baseline is a sampling-free method log likelihood maximization (LLM) (Zhang et al., 2019; Poggi et al., 2020), where a network is trained to predict a output distribution with log likelihood maximization.

C.2. Monocular Depth Estimation

For depth estimation (Postels et al., 2019; Kendall & Gal, 2017), we use one of the commonly applied models based on fully convolutional residual network (FCRN) (Laina et al., 2016). We directly use the trained model released by the original author; this is consistent with the scenarios of blackbox and graybox cases, since the code for training is not released.

We evaluate the model on NYU Depth Dataset V2. For inference, we avoid applying 90-degree rotation to input, since the orientation is a strong prior to predict depth which can violate the tolerability, and only apply horizontal flip to generate 2 samples for uncertainty estimation. For infer-dropout, we choose two locations (intermediate layers) to add the dropout layer. For infer-noise, we choose three locations to add the noise layer (two intermediate layers and one layer before the final FC layer). Then we conduct similar experiments as described in the SR task. For the baseline MC-dropout, note that the model has a dropout layer before the final fully connected (FC) layer during training, we directly perform sampling from the existing dropout layer. Sample numbers of 2 and 8 are evaluated for both infer-dropout and infer-noise.

D. Evaluation Metrics

Evaluation Metrics. Commonly used metrics to evaluate uncertainty estimation include Brier score (BS), expected calibration error (ECE), and negative log-likelihood (NLL) (Lakshminarayanan et al., 2017; Guo et al., 2017). However, BS and ECE are for classification tasks only and hence not applicable in our setting. We therefore use the following metrics for evaluations: (1) NLL, which is defined in regression tasks by assuming a Gaussian distribution. However, note that NLL depends on not only the quality of uncertainty estimation but also the prediction accuracy itself. Therefore contrary to previous belief, we argue that it is not an ideal metric for evaluating uncertainty estimation. (2) Area Under the Sparsification Error (AUSE), which quantifies how much uncertainty estimation coincides with the true errors (Ilg et al., 2018). (3) Correlation between the estimated uncertainty and the error. Here we define four variants of correlation (see details in Sec. G): pixel-wise, mean, block-wise, and patch-wise correlations to evaluate performance at the pixel, image, block, and patch levels, respectively. The intuition is that in many situations it is more instructive and meaningful when uncertainty is visualized in each region (e.g. a region with a possible tumor for a medical imaging application). Note that block-wise correlation depends on specific segmentation algorithms, while patch-wise correlation defines regions in an algorithm-independent way. Similarly we also define four evaluation forms for AUSE.

For our training-free methods, these metrics are computed between uncertainty and the error from the original model (without perturbation), because we will still use the original model for prediction. For training-required methods such as MC-dropout (i.e. MC-drop^8 in Table 3) (Gal & Ghahramani, 2016), deep ensemble (Lakshminarayanan et al., 2017) and log likelihood maximization (LLM) (Zhang et al., 2019; Poggi et al., 2020), the mean of output samples are used as prediction. Meanwhile, we also evaluate another MC-dropout variant, denoted as MC-drop^1, where the output of the original model is used as prediction, to be consistent with training-free methods.

E. Tolerability & Performance

Comparable Performance with Training-required Baselines. For the qualitative evaluation on depth estimation task shown in Fig. 4, high variance usually occurs in the area with high spatial resolution and large depth. The quantitative evaluation results using metrics described above are shown in Table 3 and Table 2. Note that SRGAN has a higher correlation than the SRresnet model. For depth estimation, we find that using infer-noise in the intermediate layers outperforms other methods. For baseline MC-dropout, we perturb right before the last convolutional layer—the only dropout layer during the original model training, and therefore produce a highly localized variance map with poor correlation, shown in Fig. 4. If we are allowed to perform MC-dropout in intermediate layers and re-train the model, a correlation value comparable to that of infer-dropout should be expected.

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Figure 4. Visualization of pixel-wise uncertainty (variance) maps from infer-transformation, infer-dropout, MC-dropout (Gal & Ghahramani, 2016) compared with the $L_1$ loss map in depth estimation task. Correlation between the $L_1$ loss map and the uncertainty map is also presented.

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Table 2. Mean/patch-wise/block-wise/pixel-wise AUSE and correlation between $L_1$ loss and uncertainty, and NLL on NYU Depth Dataset V2. Our infer-transformation, infer-dropout and infer-noise are compared with MC-dropout (Gal & Ghahramani, 2016). MC-drop$^1$ uses the output of the original model as prediction while MC-drop$^2$ uses the mean of output samples from the re-trained model (with added dropout) as prediction. Models evaluated: FCRN model.

estimation. Better tolerability means smaller decrease in accuracy after perturbation (i.e. smaller $\epsilon$ in Sec. 2.3). Fig. 5 shows the correlation for different amount of perturbations (noise or dropout) in different locations, and the corresponding predictive performance $C = |E[Z] - Y|$ (evaluated as $L_1$ loss) after perturbations. As we can see, the optimal cases to generate uncertainty maps with high correlation require that $C$ should remain small after perturbation (high tolerability).

Generally, our experiments suggest that perturbations leading to less than $20\%$ relative drop of performance work well for uncertainty estimation. Interestingly, different methods have different ways of achieving high tolerability: (1) For MC-dropout, involving dropout during training increases the robustness of model against perturbations, keeping the loss relatively small after adding dropout layer in most locations during inference; (2) for infer-dropout, adding dropout layer in intermediate locations (i.e. location 2 and location 3) where the information is the most redundant (He et al., 2014), can effectively alleviate disturbance; (3) for infer-noise, adding noise with small standard deviation effectively limits the perturbation level. More interestingly, we further find that for both MC-dropout and infer-dropout, adding perturbation in intermediate layers are usually the optimal choices for uncertainty estimation. Applying infer-dropout in these intermediate layers, we could achieve comparable or even better correlation compared to training-required baselines. For infer-noise, locations do not have similar effect; one can therefore further tune the noise strength $\sigma$ to achieve higher correlation. The conclusion above is also consistent with the evaluation of other models.

Fig. 5 shows the perturbation on the SRGAN. We plot the correlation between different amount of perturbations (noise or dropout) in different locations and the corresponding predictive performance (evaluated as $L_1$ loss) after perturbations. For SRresnet model, We find that for both MC-dropout and infer-dropout, adding perturbation in intermediate layers is usually the optimal choice for uncertainty estimation. For infer-noise, locations do not have similar effect; one can therefore further tune the noise strength $\sigma$ to
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<td>patch $L_1$</td>
<td>0.340</td>
<td>0.359</td>
</tr>
<tr>
<td>block $L_1$</td>
<td>0.501</td>
<td>0.520</td>
</tr>
<tr>
<td>pixel $L_1$</td>
<td>0.462</td>
<td>0.486</td>
</tr>
<tr>
<td>Corr</td>
<td></td>
<td></td>
</tr>
<tr>
<td>mean $L_1$</td>
<td>0.408</td>
<td>0.379</td>
</tr>
<tr>
<td>patch $L_1$</td>
<td>0.505</td>
<td>0.521</td>
</tr>
<tr>
<td>block $L_1$</td>
<td>0.237</td>
<td>0.269</td>
</tr>
<tr>
<td>pixel $L_1$</td>
<td></td>
<td></td>
</tr>
<tr>
<td>NLL</td>
<td>107,243</td>
<td>43,071</td>
</tr>
</tbody>
</table>

Table 3. Mean/patch-wise/block-wise/pixel-wise AUSE and correlation between $L_1$ loss and uncertainty, and NLL on SR benchmark dataset Set 14. Our infer-transformation, infer-dropout and infer-noise are compared with MC-dropout (Gal & Ghahramani, 2016), deep ensemble (Lakshminarayanan et al., 2017) and log likelihood maximization (LLM) (Zhang et al., 2019). MC-drop$^1$ uses the output of the original model as prediction while MC-drop$^2$ uses the mean of output samples from the re-trained model (with added dropout) as prediction. Models evaluated: SRresnet.

F. Details on Network Noise/Dropout Injection Locations

To perform uncertainty estimation using infer-noise, infer-dropout, and baseline MC-dropout on both SRGAN model and SRresnet model, we choose 4 different locations for noise injection, including the layers right after the input, as well as some intermediate layers, as shown in Fig. 7.

To perform uncertainty estimation using infer-noise and infer-dropout on FCRN model, we choose 3 different locations for noise injection, including the layers right before the output, as well as some intermediate layers, as shown in Fig. 8. For baseline MC-dropout, we choose location 3 where the dropout layer added during training for the original model.

G. Details on Evaluation Metrics

In this section, we provide the details for our proposed correlation-based evaluation metrics. Assuming we have $N$ outputs given the same input $x$ from our infer-transformation, infer-drop and infer-noise, each output is represented by $Y_w$. Given the output image with the size of $H \times W$, the error we define for regression task is pixel-wise $L_1$ loss and $L_2$ loss, represented by $L_{1,ij}$ and $L_{2,ij}$, where $i, j$ is the corresponding coordinates of the pixel $P_{ij}$ in the output image. The uncertainty (variance) estimated in these methods is also a pixel-wise value, represented by $V_{ij} = \frac{1}{N} \sum_{n=1}^{N} (Y_{w,x_{ij}} - \overline{Y}_{x_{ij}})^2$. The pixel-wise $L_1$ correlation is defined as $corr(\{V_{ij}\}, \{L_{1,ij}\})$. The second metric is mean correlation, the mean $L_1$ error $\overline{L}_{1,z} = \frac{\sum_{j=1}^{W} \sum_{i=1}^{H} L_{1,ij}}{W \times H}$ is defined as the average error of a single image $z$, correspondingly, the mean variance is defined as $\overline{V}_{z} = \frac{\sum_{j=1}^{W} \sum_{i=1}^{H} V_{ij}}{W \times H}$, the mean $L_1$ correlation is defined as $corr(\{\overline{V}_{z}\}, \{\overline{L}_{1,z}\})$. This metric has been used in (Zhang et al., 2019). The third metric for evaluation is the block-wise correlation – a new metric we propose in this work. To compute block-wise correlation, we need to firstly apply a local segmentation algorithm to the output of the trained model to cluster pixels with similar low-level context. Here we use the local-center-of-mass approach (Aganj et al., 2018) to perform segmentation. We denote each cluster as $C_i$. The variance of $K_{C_i}$ pixels inside each cluster (block) $C_i$ is then averaged and replaced with the mean value $\overline{V}_i = \frac{\sum_{C_{i} \in C} V_{ij}}{K_{C_i}}$.

The block-wise $L_1$ loss $\overline{L}_{1,i}$ can be calculated similarly. After that, we calculate the pixel-wise correlation of each pixels with the updated value as the $L_1$ block-wise correlation $corr(\{\overline{V}_{i}\}, \{\overline{L}_{1,i}\})$. For the fourth metric, patch-wise correlation, where the segmentation clusters in block-wise correlation are replaced by patches. In our analysis, each image is divided into $10 \times 10$ patches. And then the patch-wise correlation is calculated with following the same rule as block-wise correlation. Besides correlation, we also define four similar metrics in terms of AUSE. More details related with the definition of AUSE are in (Ilg et al., 2018).

Meanwhile, as illustrated in Fig. 6, we find that sparsification error has a strong association with correlation, when the oracle sparsifications of different methods are the same. As a result, when AUSE of infer-dropout and MC-dropout (defined as MC-drop$^1$ here) is nearly identical, correlation is almost the same.
**H. Details on Performance Bound**

The oracle (performance bound) is calculated with an empirically chosen constant margin $t = 5\sigma$, where $\sigma^2$ is average of $\mathbb{V}[Z]$ for all the pixels across the entire image.

**I. Applications Benefit from Uncertainty Estimation**

We find the several applications can be benefited from the uncertainty estimated in our methods. The first application is to improve the quality of SR results. We propose a novel and efficient method which takes the pixel-wise uncertainty map as a weight term for the regression loss, while keeping
Figure 7. Different locations for infer-noise and infer-dropout in SRGAN and SRresnet for super resolution. For each experiment, the noise or dropout is injected at a single location with one perturbation level.

Figure 8. Different locations for noise or dropout injection in the FCRN model for depth estimation.

<table>
<thead>
<tr>
<th>Method</th>
<th>SSIM</th>
<th>PSNR</th>
<th>$L_1$</th>
</tr>
</thead>
<tbody>
<tr>
<td>original</td>
<td>0.772</td>
<td>23.841</td>
<td>11.623</td>
</tr>
<tr>
<td>random</td>
<td>0.773</td>
<td>23.616</td>
<td>12.344</td>
</tr>
<tr>
<td>high uncertainty</td>
<td>0.790</td>
<td>23.942</td>
<td>11.497</td>
</tr>
</tbody>
</table>

Table 5. Using active learning on our generated uncertainty maps can provide a higher data efficiency. Here we select samples with high uncertainty yields better results than select randomly.

the original adversarial loss; this could provide a more photo-realistic SR output with finer structures and sharper edges. Another application is active learning (Gal et al., 2017), which aims to use uncertainty to choose next batch of data for annotation. Our result shows that active learning based on our generated uncertainty maps can provide a higher data efficiency.

The first application is to improve the quality of SR results. We propose a novel and efficient method which takes the pixel-wise uncertainty map as a weight term for the regression loss, while keeps the original adversarial loss, which could provide a more photo-realistic SR output with finer structures and sharper edges, shown in Fig. 9.

Another application is active learning (Gal et al., 2017), which aims to use uncertainty to guide annotations, then only a small subset of data are required to improve training. We find active learning based on our generated uncertainty maps can improve the performance with more efficiency, shown in Table 5.

J. Evaluations on Classification Tasks

We compare the simple training-free method using entropy and the training-required method MC-dropout on classification tasks. For classification tasks, the most straightforward and commonly used method is to calculate the entropy of output probability as uncertainty, which already qualifies as a training-free method. We then compare it with a sampling-based and training-required method – MC-dropout, tuned on different locations and using 8 samples. Here we conduct three experiments: the first one is multi-class segmentation task using Densenet (Huang et al., 2017b) on CamVid dataset; the second one is a binary segmentation task using UNET (Ronneberger et al., 2015) on a biomedical public benchmark dataset from the SNEMI3D challenge; and the third one is a classification task on CIFAR100 using ResNet (He et al., 2016). We calculate the correlation between the entropy of softmax output and the cross-entropy loss. We find that using entropy outperforms MC-dropout based on the correlation metric, as shown in Table 6.

K. Visualization of Uncertainty Maps

The uncertainty maps generated using infer-transformation, infer-dropout, infer-noise compared with MC-dropout, and
Figure 9. The first application using uncertainty map estimated in our methods to improve the quality of SR results. We compare SR results that use $L_2$ loss re-weighted by variance map (middle) and that do not (right). HR (left) represents high resolution image. Results evaluated in Set 14.

<table>
<thead>
<tr>
<th></th>
<th>Densenet</th>
<th>UNET</th>
<th>Resnet</th>
</tr>
</thead>
<tbody>
<tr>
<td>Mean Correlation</td>
<td>Entropy</td>
<td>0.928</td>
<td>0.964</td>
</tr>
<tr>
<td>Mean Correlation</td>
<td>MC-drop</td>
<td>0.718</td>
<td>0.881</td>
</tr>
<tr>
<td>Pixel-wise Correlation</td>
<td>Entropy</td>
<td>0.502</td>
<td>0.789</td>
</tr>
<tr>
<td>Pixel-wise Correlation</td>
<td>MC-drop</td>
<td>0.209</td>
<td>0.317</td>
</tr>
</tbody>
</table>

Table 6. Correlation of uncertainty and cross-entropy loss, comparing using entropy with the baseline MC-dropout, models evaluated are Densenet for the segmentation on CamVid dataset, UNET for segmentation on SNEMI3D dataset and Resnet for classification on CIFAR100 dataset.

One interesting observation is that MC-dropout tends to cap- ture local variance and ignore high-level semantics, partially because the dropout layer is always at the end of the net- work. As a result, it is difficult for MC-dropout to produce uncertainty estimates in detail-rich regions. For example, the input image in the bottom row of Fig. 11 contains a lot of details with chairs and desks. Unfortunately MC-dropout tends to ignore these details and only produce high variance in the upper half of the image (region with large depth), leading to poor correlation.
Figure 10. Visualization of uncertainty maps (log scale) and error map (log scale) from infer-transformation, infer-dropout, infer-noise compared with the baseline MC-dropout, evaluated on the SRGAN model on Set14 dataset for super-resolution task.
Figure 11. Visualization of uncertainty maps and the error map from infer-transformation, infer-dropout, infer-noise compared with the baseline MC-dropout, evaluated on the FCRN model on NYU depth dataset V2 for the depth estimation task.