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### **Bayesian Generative Active Deep Learning**

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Wed Jun 12th 04:30 - 04:35 PM @ Room 201

### Deep learning (DL)

- The dominant machine learning methodology [Huang et al., 2017, Rajkomar et al., 2018].
- Issue: significant human effort for the labeling; considerable computational resources for large-scale training process [Sun et al., 2017].

### How to address these training issues? Two popular approaches:

- (Pool-based) active learning (AL): Challenging to be applied in DL: AL may overfit the (small) informative training sets

# Main goals of this paper

 Propose a novel Bayesian generative active deep learning method

• Targets the augmentation of the labeled data set with informative generated samples



• Key technical contribution: theoretically and empirically show the informativeness of this generated sample.

Figure 1: Our proposed method depicted by VAE-ACGAN model

### Generative adversarial active learning (GAAL) [Zhu and Bento, 2017]

- Relies on an optimization problem to generate new informative samples
- Can generate rich representative training data with the assumptions:
  - GAN model has been pre-trained, and
  - The optimization during generation is solved efficiently



Figure 2: Generative adversarial active learning (GAAL) [Zhu and Bento, 2017]

### Comparison between our proposed method and GAAL [Zhu and Bento, 2017]

	GAAL [Zhu and Bento, 2017]	Ours	
acquisition function	simple (binary classifier)	more effective (deep models)	
training of the generator $(G)$	2-stage	G and $C$ are jointly trained	
and classifier $(C)$	GAN model is pre-trained	Allowing them to "co-evolve"	
classification results	not competitive enough		

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### Main technical contribution:

Combining BALD and BDA for generating new labeled samples that are informative for the training process.

- Initial labeled data:  $\mathcal{D} = \{(\mathbf{x}_i, \mathbf{y}_i)\}_{i=1}^N$ , where  $\mathbf{x}_i \in \mathcal{X} \subseteq \mathbb{R}^d$  is the data sample labeled with  $\mathbf{y}_i \in \mathcal{C} = \{1, 2, \dots, C\}$  (C = # classes).
- Bayesian active learning by disagreement (BALD) scheme [Gal et al., 2017, Houlsby et al., 2011]:
- The most informative sample x<sup>\*</sup> is selected from the (unlabeled) pool data  $\mathcal{D}_{pool}$  by [Houlsby et al., 2011]:

$$\mathbf{x}^* = \operatorname*{arg\,max}_{\mathbf{x}\in\mathcal{D}_{\mathsf{pool}}} a(\mathbf{x},\mathcal{M}),\tag{1}$$

where the acquisition function  $a(\mathbf{x}, \mathcal{M})$  is estimated by the Monte Carlo (MC) dropout method [Gal et al., 2017]

$$a(\mathbf{x}, \mathcal{M}) \approx -\sum_{c} \left(\frac{1}{T} \sum_{t} \hat{p}_{c}^{t}\right) \log \left(\frac{1}{T} \sum_{t} \hat{p}_{c}^{t}\right) + \frac{1}{T} \sum_{c,t} \hat{p}_{c}^{t} \log \hat{p}_{c}^{t}, \qquad (2)$$

where T is the number of dropout iterations,  $\hat{\mathbf{p}}^t = [\hat{p}_1^t, \dots, \hat{p}_C^t] = \operatorname{softmax}(f(\mathbf{x}; \theta^t))$ , with f is the network function parameterized by  $\theta^t \sim p(\theta|\mathcal{D})$  at the *t*-th iteration. • The generated sample x':

$$\mathbf{x}' = g(e(\mathbf{x}^*)), \tag{3}$$

where a variational autoencoder (VAE) [Kingma and Welling, 2013] contains an encoder e(.) and a decoder g(.)

• VAE training: minimizing the "reconstruction loss", where if # training iterations is sufficiently large, we have:

$$\|\mathbf{x}' - \mathbf{x}^*\| < \varepsilon, \tag{4}$$

- $\varepsilon>0$  (arbitrarily small) see Fig. 3.
- $\mathcal{D} \leftarrow \mathcal{D} \cup \{(\mathbf{x}^*, \mathbf{y}^*), (\mathbf{x}^{'}, \mathbf{y}^*)\}$ , which are used for the next training iteration.



**Figure 3:** Reduction of  $||\mathbf{x}' - \mathbf{x}^*||$  as the training progresses of the VAE model.

## **Key Question**

How about the "information content" of the generated sample  $\mathbf{x}'$ measured by  $a(\mathbf{x}', \mathcal{M})$ ?

### **Proposition 2.1**

Assuming that there exists the gradient of the acquisition function  $a(\mathbf{x}, \mathcal{M})$  with respect to the variable  $\mathbf{x}$ , namely  $\nabla_{\mathbf{x}} a(\mathbf{x}, \mathcal{M})$ , and that  $\mathbf{x}^*$  is an interior point of  $\mathcal{D}_{pool}$ , then  $a(\mathbf{x}', \mathcal{M}) \approx a(\mathbf{x}^*, \mathcal{M})$  (i.e., the absolute difference between these values are within a certain range). Consequently, the sample  $\mathbf{x}'$  generated from the most informative sample  $\mathbf{x}^*$  by (3) is also informative.

![](_page_7_Figure_1.jpeg)

Figure 4: Network architecture of our proposed model.

- Classification performance measured by the top-1 accuracy as a function of the number of acquisition iterations and the percentage of training samples.
- Our proposed algorithm, active learning using "information-preserving" data augmentation (AL w. VAEACGAN) is compared with:
  - Active learning using BDA (AL w. ACGAN)
  - BALD [Gal et al., 2017] without using data augmentation (AL without DA),
  - BDA [Tran et al., 2017] without active learning (BDA) (using full and partial training sets)
  - Random selection
- Benchmark data sets: MNIST [LeCun et al., 1998], CIFAR-10, CIFAR-100 [Krizhevsky et al., 2012], and SVHN [Netzer et al., 2011].
- Baseline classifiers: ResNet18 [He et al., 2016a] and ResNet18pa [He et al., 2016b]

![](_page_9_Figure_1.jpeg)

**Figure 5:** Training and classification performance of the proposed Bayesian generative active learning (AL w. VAEACGAN) compared to other methods. This performance is measured as a function of the number of acquisition iterations and respective percentage of samples from the original training set used for modeling.

Table I: Mean  $\pm$  standard deviation of the classification accuracy on MNIST, CIFAR-10, and CIFAR-100 after 150 iterations over 3 runs

			MNIST				
	AL W. VAEACGAN	AL W. ACGAN	AL W. PMDA	AL WITHOUT DA	BDA (partial training)	RANDOM SELECTION	
Resnet18	$99.53 \pm 0.05$	$99.45 \pm 0.02$	$99.37 \pm 0.15$	$99.33 \pm 0.10$	$99.33 \pm 0.04$	$99.00 \pm 0.13$	
Resnet18pa	$99.68 \pm 0.08$	$99.57 \pm 0.07$	$99.49 \pm 0.09$	$99.35 \pm 0.11$	$99.35 \pm 0.07$	$99.20 \pm 0.12$	
CIFAR-10							
Resnet18	$87.63 \pm 0.11$	$86.80 \pm 0.45$	$82.17 \pm 0.35$	$79.72 \pm 0.19$	$85.08 \pm 0.31$	$77.29 \pm 0.23$	
Resnet18pa	$91.13 \pm 0.10$	$90.70 \pm 0.24$	$87.70 \pm 0.39$	$85.51 \pm 0.21$	$86.90 \pm 0.27$	$80.69 \pm 0.19$	
CIFAR-100							
Resnet18	$68.05 \pm 0.17$	$66.50 \pm 0.63$	$55.24 \pm 0.57$	$50.57 \pm 0.20$	$65.76 \pm 0.40$	$49.67 \pm 0.52$	
Resnet18pa	$69.69 \pm 0.13$	$67.79 \pm 0.76$	$59.67 \pm 0.60$	$55.82 \pm 0.31$	$65.79 \pm 0.51$	$54.77 \pm 0.29$	

![](_page_10_Picture_3.jpeg)

Figure 6: Images generated by our proposed (AL w. VAEACGAN) approach for each data set.

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- Consistently shows that our proposed approach outperforms other methods.
- The classification results by AL w. VAEACGAN are statistically significant w.r.t. BDA (partial training) on all those data sets, and w.r.t. AL w. ACGAN on CIFAR- $\{10, 100\}$  for both models (i.e.,  $p \le .05$ , two-sample t-test).
- Proposed a novel Bayesian generative active deep learning approach that consistently shows to be more effective than data augmentation and active learning in several classification problems.
- Our proposed approach is (active learning) model-agnostic, it therefore can be combined with several currently sate-of-the-art active learning methods [Ducoffe and Precioso, 2018, Gissin and Shalev-Shwartz, 2018, Sener and Savarese, 2018].

### **Future work**

- To investigate how to generate samples directly using complex acquisition functions
- To work on the efficiency of our proposed method (its empirical computational cost is slightly higher than BDA [Tran et al., 2017] and BALD [Gal et al., 2017, Houlsby et al., 2011]).
- Imbalanced data sets

# Poster presentation: Wed Jun 12th 06:30–09:00 PM @ Pacific Ballroom #263

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