Dichotomize and Generalize: PAC-Bayesian Binary Activated Deep Neural Networks



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Introduction

We present a comprehensive study of multilayer neural networks with binary activation, relying on the PAC-Bayesian theory.

Contributions

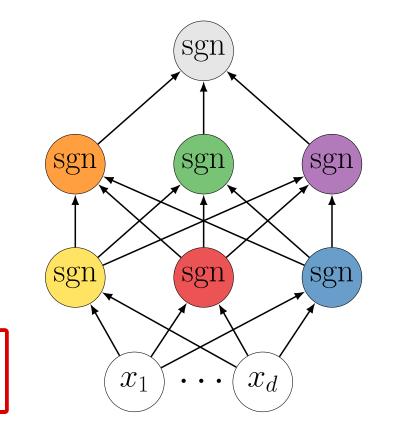
- ► An end-to-end framework to train a binary activated deep neural network (DNN).
- ► Nonvacuous PAC-Bayesian generalization bounds for binary activated DNNs.

Binary Activated Neural Networks

- ightharpoonup L fully connected layers
- $ightharpoonup d_k$ denotes the number of neurons of the $k^{
 m th}$ layer
- $ightharpoonup \operatorname{sgn}(a) = 1 \text{ if } a > 0 \text{ and } \operatorname{sgn}(a) = -1 \text{ otherwise}$
- $lackbox{Weights matrices}: \mathbf{W}_k \in \mathbb{R}^{d_k \times d_{k-1}}, \ \theta = \mathrm{vec} \big(\{\mathbf{W}_k\}_{k=1}^L \big) \in \mathbb{R}^D$



 $f_{\theta}(\mathbf{x}) = \operatorname{sgn}(\mathbf{w}_{L}\operatorname{sgn}(\mathbf{W}_{L-1}\operatorname{sgn}(\ldots\operatorname{sgn}(\mathbf{W}_{1}\mathbf{x}))))$



PAC-Bayesian Theory

Given a data distribution \mathcal{D} , a training set $S = \{(\mathbf{x}_1, y_1), \dots, (\mathbf{x}_n, y_n)\} \sim \mathcal{D}^n$, with $\mathbf{x}_i \in \mathbb{R}^{d_0}$ and $y_i \in \{-1, 1\}$, a loss $\ell : [-1, 1]^2 \to [0, 1]$, a predictor $f \in \mathcal{F}$:

$$\mathcal{L}_{\mathcal{D}}(f) = \underset{(\mathbf{x},y)\sim\mathcal{D}}{\mathbf{E}} \; \ell\Big(f(\mathbf{x}),y\Big) \qquad \longleftarrow \underline{\text{generalization loss}}$$

$$\widehat{\mathcal{L}}_S(f) = rac{1}{n} \sum_{i=1}^n \ell \Big(f(\mathbf{x}_i), y_i \Big) \qquad \longleftarrow \underbrace{\textit{empirical loss}}$$

— PAC-Bayesian Theorem -

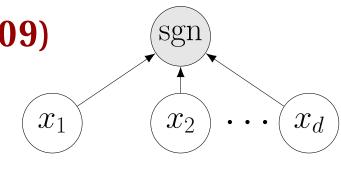
For any prior P on \mathcal{F} , with probability $1-\delta$ on the choice of $S\sim\mathcal{D}^n$, we have for all C > 0, and any posterior distribution Q on \mathcal{F} :

$$\underset{f \sim Q}{\mathbf{E}} \mathcal{L}_{\mathcal{D}}(f) \leq \frac{1}{1 - e^{-C}} \left(1 - \exp\left(-C \underset{f \sim Q}{\mathbf{E}} \widehat{\mathcal{L}}_{S}(f) - \frac{1}{n} [\mathrm{KL}(Q \| P) + \ln \frac{2\sqrt{n}}{\delta}] \right) \right)$$

Linear Classifier

PAC-Bayesian Learning of Linear Classifiers (Germain et al., 2009)

$$f_{\mathbf{w}}(\mathbf{x}) \coloneqq \operatorname{sgn}(\mathbf{w} \cdot \mathbf{x})$$
, with $\mathbf{w} \in \mathbb{R}^d$



PAC-Bayesian analysis

- ▶ Space of all linear classifiers $\mathcal{F}_d \coloneqq \{f_{\mathbf{v}} | \mathbf{v} \in \mathbb{R}^d\}$
- ► Gaussian prior $P_{\mathbf{u}} \coloneqq \mathcal{N}(\mathbf{u}, I_d)$ over \mathcal{F}_d
- ► Gaussian posterior $Q_{\mathbf{w}} \coloneqq \mathcal{N}(\mathbf{w}, I_d)$ over \mathcal{F}_d
- ▶ Predictor $F_{\mathbf{w}}(\mathbf{x}) \coloneqq \mathbf{E}_{\mathbf{v} \sim Q_{\mathbf{w}}} f_{\mathbf{v}}(\mathbf{x}) = \operatorname{erf}\left(\frac{\mathbf{w} \cdot \mathbf{x}}{\sqrt{2}\|\mathbf{x}\|}\right)$
- Linear loss $\ell(f_{\mathbf{v}}(\mathbf{x}), y) \coloneqq \frac{1}{2} \frac{1}{2}yf_{\mathbf{v}}(\mathbf{x})$

tanh(x)

Bound minimization

$$C n \widehat{\mathcal{L}}_S(F_{\mathbf{w}}) + \text{KL}(Q_{\mathbf{w}} || P_{\mathbf{u}}) = C \frac{n}{2} + C \frac{1}{2} \sum_{i=1}^n \text{erf}\left(-y_i \frac{\mathbf{w} \cdot \mathbf{x}_i}{\sqrt{2} || \mathbf{x}_i ||}\right) + \frac{1}{2} || \mathbf{w} - \mathbf{u} ||^2$$

Shallow Learning

Posterior $Q_{\theta} = \mathcal{N}(\theta, I_D)$, over the family of all networks $\mathcal{F}_D = \{f_{\tilde{\theta}} \mid \theta \in \mathbb{R}^D\}$, where $f_{\theta}(\mathbf{x}) = \operatorname{sgn}(\mathbf{w}_2 \cdot \operatorname{sgn}(\mathbf{W}_1 \mathbf{x}))$

$$F_{\theta}(\mathbf{x}) = \sum_{\tilde{\theta} \sim Q_{\theta}} f_{\tilde{\theta}}(\mathbf{x})$$

$$= \int_{\mathbb{R}^{d_{1} \times d_{0}}} Q_{1}(\mathbf{V}_{1}) \int_{\mathbb{R}^{d_{1}}} Q_{2}(\mathbf{v}_{2}) \operatorname{sgn}(\mathbf{v}_{2} \cdot \operatorname{sgn}(\mathbf{V}_{1}\mathbf{x})) d\mathbf{v}_{2} d\mathbf{V}_{1}$$

$$= \sum_{\mathbf{s} \in \{-1,1\}^{d_{1}}} \operatorname{erf}\left(\frac{\mathbf{w}_{2} \cdot \mathbf{s}}{\sqrt{2d_{1}}}\right) \int_{\mathbb{R}^{d_{1} \times d_{0}}} \mathbf{1} [\mathbf{s} = \operatorname{sgn}(\mathbf{V}_{1}\mathbf{x})] Q_{1}(\mathbf{V}_{1}) d\mathbf{V}_{1}$$

$$= \sum_{\mathbf{s} \in \{-1,1\}^{d_{1}}} \operatorname{erf}\left(\frac{\mathbf{w}_{2} \cdot \mathbf{s}}{\sqrt{2d_{1}}}\right) \underbrace{\prod_{i=1}^{d_{1}} \left[\frac{1}{2} + \frac{s_{i}}{2} \operatorname{erf}\left(\frac{\mathbf{w}_{1}^{i} \cdot \mathbf{x}}{\sqrt{2} \|\mathbf{x}\|}\right)\right]}_{Pr(\mathbf{s}|\mathbf{x},\mathbf{W}_{1})}$$

PAC-Bayesian bound ingredients

- ► Empirical loss : $\widehat{\mathcal{L}}_S(F_{\theta}) = \mathbf{E}_{\theta' \sim Q_{\theta}} \widehat{\mathcal{L}}_S(f_{\theta'}) = \frac{1}{n} \sum_{i=1}^n \left[\frac{1}{2} \frac{1}{2} y_i F_{\theta}(\mathbf{x}_i) \right]$
- ► Complexity term : $\mathrm{KL}(Q_{\theta} \| P_{\mu}) = \frac{1}{2} \| \theta \mu \|^2$, with $\mu \in \mathbb{R}^D$
- ► Generalization bound: $\frac{1}{1-e^{-C}} \left(1 \exp\left(-C \widehat{\mathcal{L}}_S(F_\theta) \frac{1}{n} [\text{KL}(Q_\theta || P_\mu) + \ln \frac{2\sqrt{n}}{\delta}] \right) \right)$

Contact: gael.letarte.l@ulaval.ca Code: github.com/gletarte/dichotomize-and-generalize



Visualization

The proposed method can be interpreted as a majority vote of hidden layer representations.

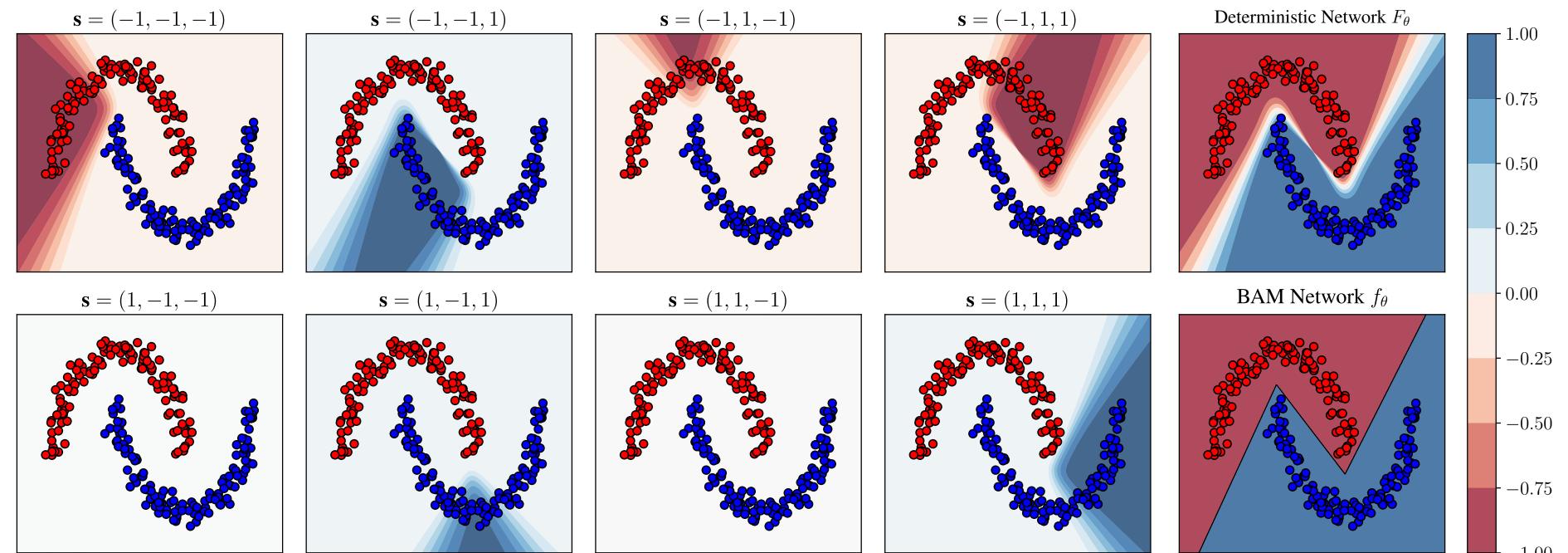


FIGURE 1: Illustration of the proposed method for a one hidden layer network of size $d_1=3$, interpreted as a majority vote over 8 binary representations $\mathbf{s} \in \{-1, 1\}^3$. For each \mathbf{s} , a plot shows the values of $F_{\mathbf{w}_2}(\mathbf{s}) \Pr(\mathbf{s} | \mathbf{x}, \mathbf{W}_1)$.

Stochastic Approximation

Prediction.

Hidden layer partial derivatives, with $erf'(x) := \frac{2}{\sqrt{\pi}}e^{-x^2}$.

$$F_{\theta}(\mathbf{x}) = \sum_{\mathbf{s} \in \{-1,1\}^{d_1}} F_{\mathbf{w}_2}(\mathbf{s}) \Pr(\mathbf{s} | \mathbf{x}, \mathbf{W}_1)$$

$$\frac{\partial}{\partial \mathbf{w}_{1}^{k}} F_{\theta}(\mathbf{x}) = \frac{\mathbf{x}}{2^{\frac{3}{2}} \|\mathbf{x}\|} \operatorname{erf}' \left(\frac{\mathbf{w}_{1}^{k} \cdot \mathbf{x}}{\sqrt{2} \|\mathbf{x}\|} \right) \sum_{\mathbf{s} \in \{-1,1\}^{d_{1}}} s_{k} F_{\mathbf{w}_{2}}(\mathbf{s}) \left[\frac{\operatorname{Pr}(\mathbf{s}|\mathbf{x}, \mathbf{W}_{1})}{\operatorname{Pr}(s_{k}|\mathbf{x}, \mathbf{w}_{1}^{k})} \right]$$

— Monte Carlo sampling — We generate T random binary vectors $\{\mathbf{s}^t\}_{t=1}^T$ according to $\Pr(\mathbf{s}|\mathbf{x},\mathbf{W}_1)$.

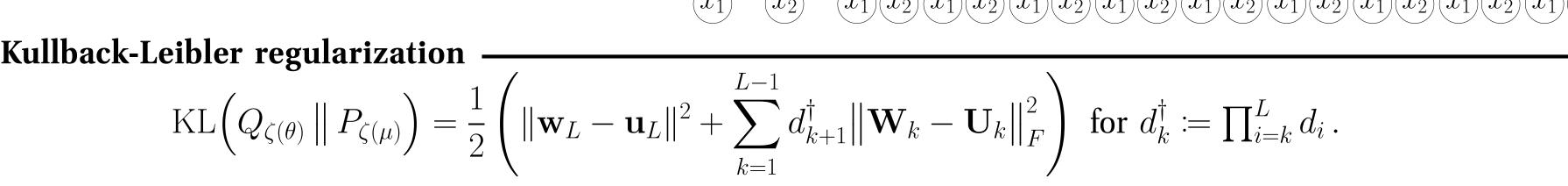
$$F_{\theta}(\mathbf{x}) \approx \frac{1}{T} \sum_{t=1}^{T} F_{\mathbf{w}_2}(\mathbf{s}^t) \qquad \qquad \frac{\partial}{\partial \mathbf{w}_1^k} F_{\theta}(\mathbf{x}) \approx \frac{\mathbf{x}}{2^{\frac{3}{2}} \|\mathbf{x}\|} \operatorname{erf}'\left(\frac{\mathbf{w}_1^k \cdot \mathbf{x}}{\sqrt{2} \|\mathbf{x}\|}\right) \frac{1}{T} \sum_{t=1}^{T} \frac{s_k^t}{\Pr(s_k^t | \mathbf{x}, \mathbf{w}_1^k)} F_{\mathbf{w}_2}(\mathbf{s}^t)$$

This turns out to be a variant of the REINFORCE algorithm (Williams, 1992).

Deep Learning (PBGNet)

To enable a layer-by-layer computation of the prediction function, we want the neurons of a given layer to be independent of each other. This is achieved with the tree architecture mapping function $\zeta(\theta)$ applied on a multilayer network.

Recursive definition. $F_k^{(j)}$ denotes the output of the j^{th} neuron of the k^{th} hidden layer : $F_1^{(j)}(\mathbf{x}) = \operatorname{erf}\left(\frac{\mathbf{w}_1^j \cdot \mathbf{x}}{\sqrt{2}\|\mathbf{x}\|}\right)$ first hidden layer $F_{k+1}^{(j)}(\mathbf{x}) = \sum_{\mathbf{s} \in \{-1,1\}^{d_k}} \operatorname{erf}\left(\frac{\mathbf{w}_{k+1}^{j} \cdot \mathbf{s}}{\sqrt{2d_k}}\right) \prod_{i=1}^{d_k} \left(\frac{1}{2} + \frac{1}{2}s_i \times F_k^{(i)}(\mathbf{x})\right)$



Experiment

We perform model selection using a validation set for a MLP with tanh activations, and using the PAC-Bayes bound for our PBGnet and PBGnet_{pre} algorithms. PBGnet_ℓ and PBGnet_{ℓ-bnd} are intermediate variants.

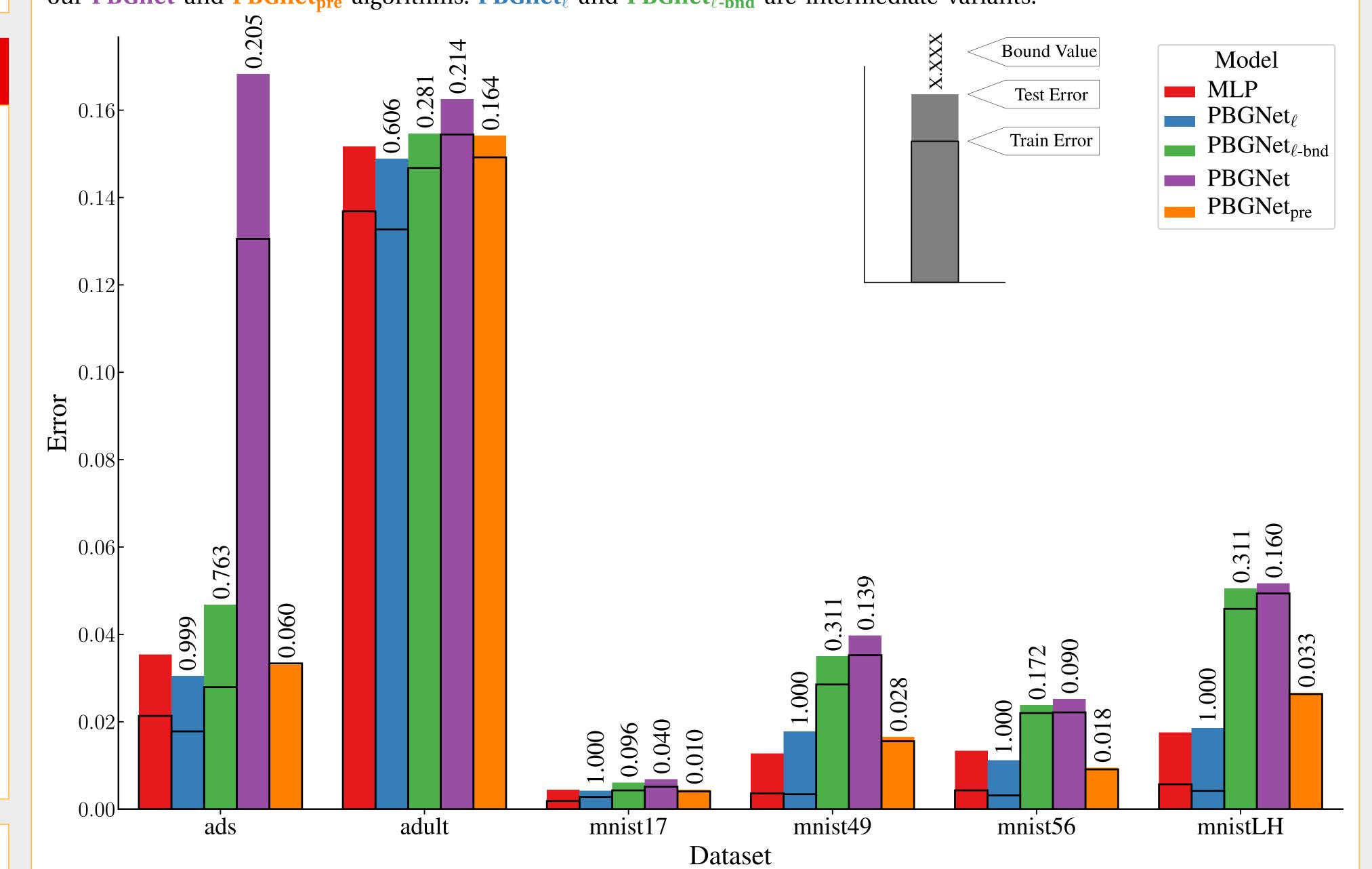


FIGURE 2: Experiment results for the considered models on the binary classification datasets. PAC-Bayesian bounds hold with probability 0.95.