On Uncertainty Estimation by Tree-based Surrogate Models in Sequential Model-based Optimization

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Introduction

- In sequential model-based optimization, GP regression [Rasmussen and Williams, 2006] is a popular choice as a surrogate model, because of its capability of calculating uncertainties analytically.
- On the other hand, an ensemble of randomized trees is another option and has practical merits over GPs due to its scalability and easiness of handling continuous/discrete mixed variables [Hutter et al., 2011].
- We revisit various ensembles of randomized trees to investigate their behavior in the perspective of prediction uncertainty estimation.
- Then, we propose our own tree-based model, referred to as *BwO forest*.

Uncertainty Estimation by Tree-based Models

- **Sum-of-Trees Model**: Posterior mean and variance functions are

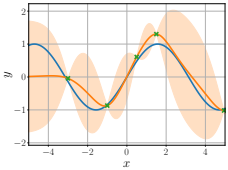
$$\mu\big(\mathbf{x}; \{\mathcal{T}_b\}_{b=1}^{B}, \mathbf{X}, \mathbf{y}\big) = \frac{1}{B}\sum_{b=1}^{B}\sum_{\tau\in\boldsymbol{\tau}_{b,l}} \mu_\tau 1_{\mathbf{x}\in\tau}, \tag{1}$$

$$\sigma^2\big(\mathbf{x}; \{\mathcal{T}_b\}_{b=1}^{B}, \mathbf{X}, \mathbf{y}\big) = \frac{1}{B}\sum_{b=1}^{B}\Bigg(\Big(\sum_{\tau\in\boldsymbol{\tau}_{b,l}} \sigma_\tau 1_{\mathbf{x}\in\tau}\Big)^2 + \Big(\sum_{\tau\in\boldsymbol{\tau}_{b,l}} \mu_\tau 1_{\mathbf{x}\in\tau}\Big)^2\Bigg) - \mu\big(\mathbf{x}; \{\mathcal{T}_b\}_{b=1}^{B}, \mathbf{X}, \mathbf{y}\big)^2, \tag{2}$$

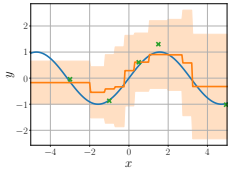
  by the law of total variance, as described in [Hutter et al., 2014].

- **Gradient Boosting Models**: It updates parameters $\boldsymbol{\theta}$ using their gradients in terms of the objective of parametric dist., e.g., likelihood function.

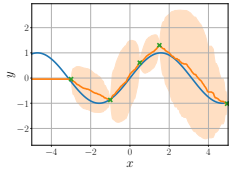
Uncertainty Estimation by Tree-based Models



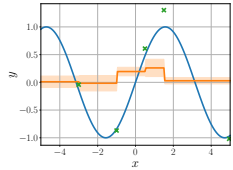
(a) Gaussian process



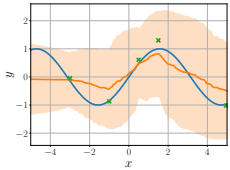
(b) Random forest



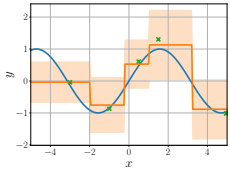
(c) Extremely randomized trees



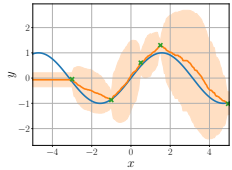
(d) BART



(e) Mondrian forest



(f) NGBoost



(g) BwO forest

Figure 1: Results with GP regression and tree-based models such as random forest, extremely randomized trees, BART, Mondrian forest, NGBoost, and BwO forest (ours).

Tree Construction

- The uncertainty of an ensemble is derived from the randomness of individual trees:
  - (i) **bagging** [Breiman, 1996]: it samples a bootstrap sample with replacement and then aggregates base estimators;
  - (ii) **random feature selection**: this technique randomly selects a feature from a set of dimensions;
  - (iii) **random split location**: it randomly selects a split location between lower and upper bounds of the selected dimension;
  - (iv) **random tree sampling**: this strategy randomly samples a tree under the assumption on a prior distribution.

- Random forest [Breiman, 2001] employs (i) and (ii), extremely randomized trees [Geurts et al., 2006] employs (ii) and (iii), BART [Chipman et al., 2010] employs (i), (ii), and (iv), and Mondrian forest [Lakshminarayanan et al., 2016] employs (i) and (iii).

- On the contrary, NGBoost [Duan et al., 2020] is defined as a gradient boosting model.

Elaborating Uncertainty Estimation by Tree-based Models

- As pointed out in the work [Mendelson et al., 2016], the expectation and variance of an indicator for the existence of $\mathbf{x}_i$ in a bootstrap sample $\mathbf{B}$ are expressed as

$$\mathbb{E}[1_{\mathbf{x}_i \in \mathbf{B}}] = 1 - \left(1 - \frac{1}{N}\right)^M, \tag{3}$$

$$\mathrm{Var}[1_{\mathbf{x}_i \in \mathbf{B}}] = \left(1 - \frac{1}{N}\right)^M - \left(1 - \frac{1}{N}\right)^{2M}, \tag{4}$$

  where N is the size of $\mathbf{X}$ and M is the size of a bootstrap sample.

- The distribution of unique original elements in a bootstrap sample $\mathbf{B}$ is:

$$\mathbb{E}[|\mathrm{unique}(\mathbf{B})|] = N - \frac{(N-1)^M}{N^{M-1}}, \tag{5}$$

$$\mathrm{Var}[|\mathrm{unique}(\mathbf{B})|] = (N-1)\frac{(N-2)^M}{N^{M-1}} + \frac{(N-1)^M}{N^{M-1}} - \frac{(N-1)^{2M}}{N^{2M-2}}, \tag{6}$$

  where $\mathrm{unique}(\mathbf{B})$ filters duplicates.

Elaborating Uncertainty Estimation by Tree-based Models

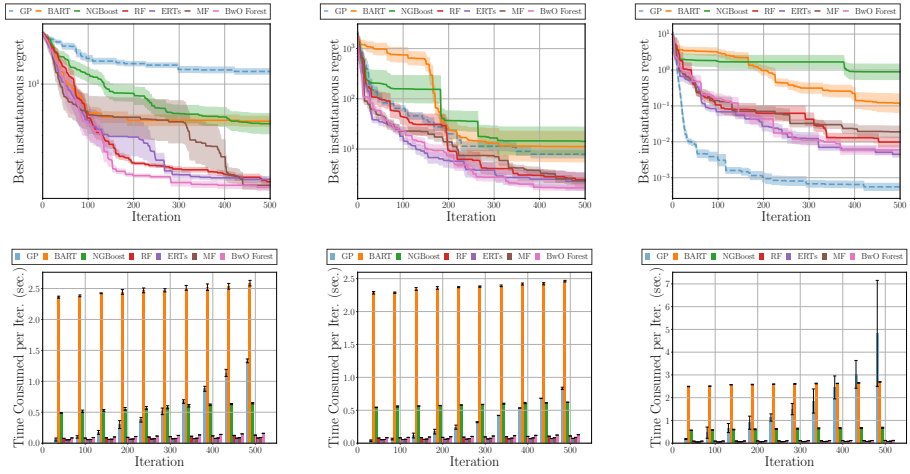
Algorithm 1 Training BwO Forest

Input: Ensemble size B, training data and evaluations $\mathbf{X} \in \mathbb{R}^{N \times d}$ and $\mathbf{y} \in \mathbb{R}^{N}$, size of bootstrap sample $M = \alpha N$ for an oversampling rate $\alpha > 1$.

Output: Set of decision trees $\{\mathcal{T}_b\}_{b=1}^{B}$

1: Initialize a set of decision trees.
2: **for** $b = 1, \ldots, B$ **do**
3: Sample a bootstrap sample $\mathbf{B}_b \in \mathbb{R}^{M \times d}$ from $\mathbf{X}$.
4: Set a root node $\boldsymbol{\tau}_r$ that contains all the elements in $\mathbf{B}_b$, and $\boldsymbol{\tau}_r$ is set as the current split node.
5: Train a decision tree using random feature selection and random split location.
6: **end for**
7: **return** A set of decision trees $\{\mathcal{T}_b\}_{b=1}^{B}$

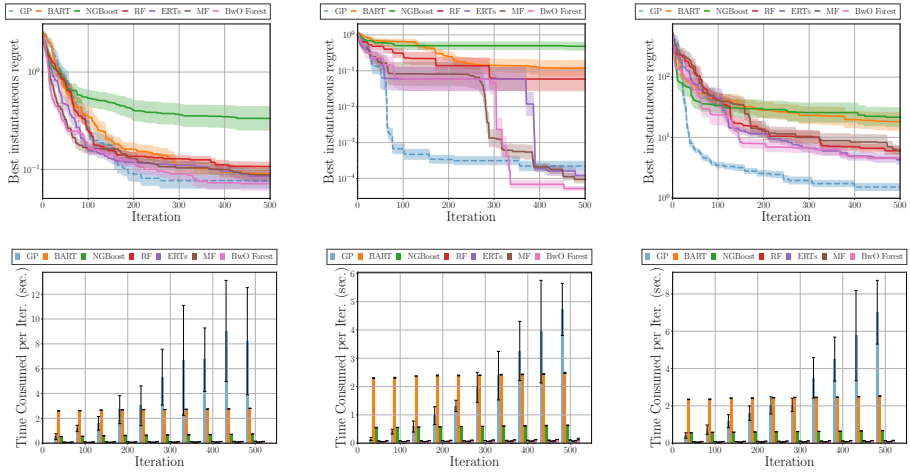
Experimental Results



(a) Ackley (4D) (b) Bohachevsky (c) Branin

Figure 2: Experimental results on various benchmark functions.

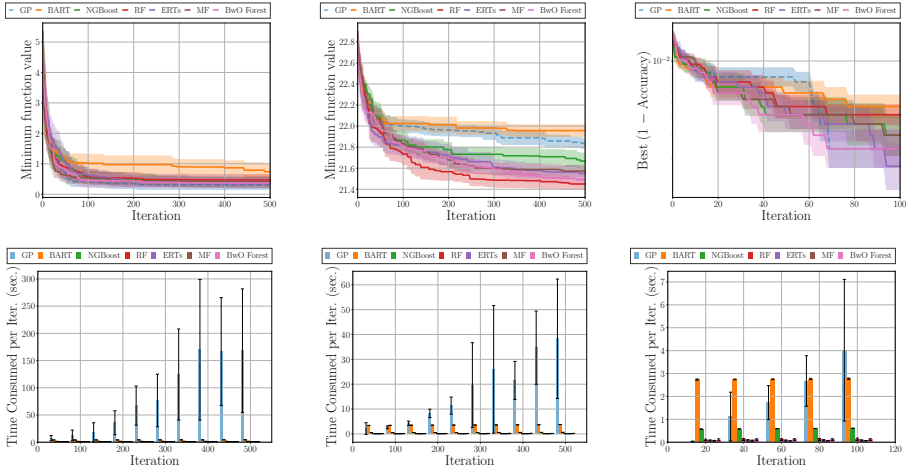
Experimental Results



(a) Hartmann6D (b) Michalewicz (c) Rosenbrock (4D)

Figure 3: Experimental results on various benchmark functions.

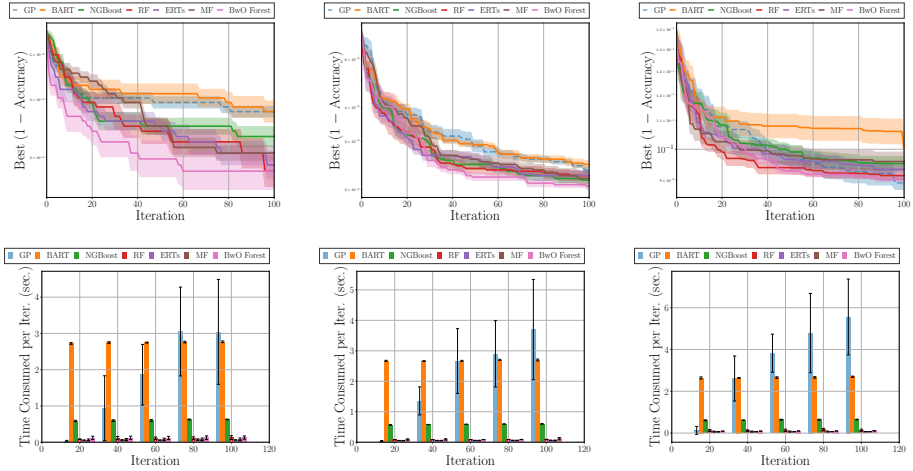
Experimental Results



(a) Ising (24D, $\lambda = 0.01$) (b) Cont. (25D, $\lambda = 0.01$) (c) Authorship

Figure 4: Experimental results on two functions defined on high-dimensional binary search spaces and automated machine learning for the Authorship dataset.

Experimental Results



(a) Breast Cancer (b) Digits (c) Phoneme

Figure 5: Experimental results on automated machine learning for three datasets, Breast Cancer, Digits, and Phoneme.

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Thank you!

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