# **Uncertainty-aware Accumulated Local Effects (UALE) for quantifying the heterogeneity of instance-level feature effects**

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# **Abstract**

Accumulated Local Effects (ALE) is a popular explainable AI method that quantifies how a feature influences the decisions of a model, handling well feature correlations. In case of complex interactions between features, instancelevel feature effects may deviate from the ALE curve. It is therefore crucial to quantify this deviation, namely, the uncertainty of the effect. In this work, we define Uncertainty-aware ALE (UALE) to quantify and visualize, on a single plot, both the average effect and its uncertainty. We show that UALE quantifies uncertainty effectively, even in case of correlated features. We also note that as in ALE, UALE's approximation requires partitioning the feature domain into nonoverlapping intervals (bin-splitting). The average effect and the uncertainty are computed from the instances that lie in each bin. We formally prove that to achieve an unbiased approximation of the uncertainty in each bin, bin-splitting must follow specific constraints. Based on this, we propose a method to determine the optimal intervals, [balancing the estimation bias and variance.] We demonstrate, through synthetic and real datasets, (a) the advantages of modeling the uncertainty with UALE compared to alternative methods and (b) the effectiveness of UALE's appropriate bin splitting for a [good] approximating of the average effect and its uncertainty.

# 1 INTRODUCTION

Recently, Machine Learning (ML) has been widely adopted in mission critical domains, such as healthcare and finance. In such high-stakes areas, it is important not only to proFor this reason, there is an increased interest in Explainable AI (XAI). XAI literature distinguishes between local and global methods (Molnar et al., 2020a). Local methods provide instance-level explanations, i.e., explain the prediction for a specific instance, whereas global methods summarize the entire model behavior. Global methods create a global explanation by aggregating the instance-level explanations into a single interpretable outcome, usually a number or a plot.

vide accurate predictions but also meaningful explanations.

A popular class of global XAI are feature effect (FE) methods (Grömping, 2020) that quantify the average (over all instances) effect of a single feature on the output<sup>1</sup>. There are three widely-used FE methods: *Partial Dependence Plots* (PDP)(Friedman, 2001), *Marginal Plots* (MP)(Apley and Zhu, 2020) and *Accumulated Local Effects* (ALE)(Apley and Zhu, 2020). ALE is established as the state-of-the-art in quantifying the average effect, since PDP and MP have been criticized (Grömping, 2020) of providing misleading explanations in case of correlated input features.

When complex interactions between features exist, the instance-level (local) feature effects may exhibit significant deviation, aka heterogeneity, from the aggregated (global) explanation, a phenomenon known as aggregation bias (Mehrabi et al., 2021). We call this deviation the uncertainty of the global effect. Aggregation bias leads to ambiguities; for example, a feature with zero average effect may indicate (a) no effect on the output or (b) an effect that is highly positive for some instances and highly negative for some others. As a result, there is a need for FE methods to quantify the deviation (uncertainty) of instance-level explanations, in addition to the average effect. In the case of PDP, such deviation is visualized via the Individual Conditional Explanations (ICE) (Goldstein et al., 2015). ICE plots, however, have the same limitations as PDP in case of correlated features. No method to quantify the deviation of instance-level effects has been proposed for ALE.

In this work, we propose UALE, a global feature-effect

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<sup>&</sup>lt;sup>1</sup>FE methods also isolate the combined effect of a pair of features to the output. Combinations of more than two features are uncommon, since they are difficult to estimate and visualize.

method that extends ALE for modeling the uncertainty of the feature effect. Our method handles well cases with correlated features. As in ALE, we provide an interval-based formulation of UALE. This formulation partitions the feature domain into non-overlapping intervals (bin-splitting) and defines the average effect (bin-effect) and its uncertainty (bin-uncertainty). We exploit this formulation to provide an UALE approximation. As we formally prove, a fixed equi-width partitioning that does not consider the characteristics of the instance-level effects, which is the case in ALE, may lead to an erroneous (biased) estimation of the uncertainty. Therefore, we propose a variable-width partitioning that leads to an unbiased approximation.

The contributions of this paper are:

- A global feature effect method (UALE) that quantifies both the average effect and its uncertainty, i.e., the deviation of instance-level effects.
- An unbiased approximation of the uncertainty through a dynamic variable-size bin partitioning of the feature domain.
- The evaluation of UALE on both synthetic and real datasets.

The implementation of our method and the code for reproducing all the experiments is provided along with the manuscript and will become publicly available upon acceptance.

# 2 BACKGROUND AND RELATED WORK

In the rest of the paper, we use following notation. Let  $\mathcal{X} \in \mathbb{R}^d$  be the d-dimensional feature space,  $\mathcal{Y} \in \mathbb{R}$  the target space and  $f(\cdot): \mathcal{X} \to \mathcal{Y}$  the black-box function. We use index  $s \in \{1,\ldots,d\}$  for the feature of interest and  $c = \{1,\ldots,d\}-s$  for the set with all other indexes. For convenience, we denote the feature vector  $\mathbf{x} = (x_1,\cdots,x_s,\cdots,x_D)$  with  $(x_s,\mathbf{x}_c)$  and the corresponding random variables  $X = (X_1,\cdots,X_s,\cdots,X_D)$  with  $(X_s,X_c)$ . The training set is  $\mathcal{D} = \{(\mathbf{x}^i,y^i)\}_{i=1}^N$  sampled i.i.d. from the distribution  $\mathbb{P}_{X,Y}$ . Finally, we use  $f^{<\mathtt{method}>}(x_s)$  for denoting the s-th FE, where  $<\mathtt{method}>$  indicates the particular method in use, for example ALE. An extensive list with all symbols used in this paper is provided at Appendix A.

#### 2.1 Feature Effect Methods

The three well-known feature effect methods are:  $Partial\ Dependence\ Plots\ (PDP),\ Marginal\ Plots\ (MP)$  and  $Accumulated\ Local\ Effects\ (ALE).\ PDP\ formulates\ the$  global FE as an expectation over the distribution of  $X_c$ ,

i.e.,  $f^{\text{PDP}}(x_s) = \mathbb{E}_{X_c}[f(x_s, X_c)]$ , whereas MP as an expectation over the distribution of  $X_c|X_s$ , i.e.,  $f^{\text{MP}}(x_s) = \mathbb{E}_{X_c|x_s}[f(x_s, X_c)]$ . Both methods suffer from misestimations in case of correlated features. PDP integrates over unrealistic instances and MP computes aggregated effects, i.e., imputes the combined effect of sets of features to a single feature (Apley and Zhu, 2020).

ALE overcomes these problems. Specifically, ALE defines the local effect of the s-th feature at a specific point  $(z, \mathbf{x}_c)$  of the input space  $\mathcal{X}$  with the partial derivative  $f^s(z, \mathbf{x}_c) = \frac{\partial f}{\partial x_s}(z, \mathbf{x}_c)$ . All the local explanations at z are then weighted by the conditional distribution  $p(\mathbf{x}_c|z)$  and are averaged, to produce the averaged effect  $\mu(z)$ . ALE plot is the accumulation of the averaged local effects:

$$f^{\text{ALE}}(x_s) = \int_{x_{s,min}}^{x_s} \underbrace{\mathbb{E}_{X_c|X_s=z}\left[f^s(z, X_c)\right]}_{u(z)} \partial z \qquad (1)$$

where  $x_{s,min}$  is the minimum value of the s-th feature.

#### 2.2 Heterogeneity Of Instance-Level Effects

The global effect is computed as an expectation over local (instance-level) effects. In addition to this global effect, it is important to know to what extent the local effects deviate from the global explanation.

ICE plots and similar methods (e.g., d-ICE plots (Goldstein et al., 2015)) provide a set of curves illustrated on top-of PDP. Each curve corresponds to one instance of the dataset,  $f_i^{\rm ICE}(x_s) = f(x_s, \mathbf{x}_c^i)$ . The user then visually observes if the curves are homogeneous, i.e., all instances have similar effect plots, and to what extent they deviate from the PDP. There are methods try to automate the aforementioned visual exploration, by grouping ICE plots into clusters (Molnar et al., 2020b; Herbinger et al., 2022; Britton, 2019). Unfortunately, these methods are subject to the failure modes of PDPs in cases of correlated features (Baniecki et al., 2021), as we also confirm in our experimental analysis (c.f., Section 4.1).

Other approaches, like H-Statistic (Friedman and Popescu, 2008), Greenwell's interaction index (Greenwell et al., 2018) or SHAP interaction values (Lundberg et al., 2018) quantify interactions between the input features. A strong interaction is an implicit indication of the existence of heterogeneous effects. These methods, however, do not quantify directly the level of heterogeneity.

To the best of our knowledge, no work exist so far that quantifies the heterogeneous effects (uncertainty) based on the formulation of ALE.

#### 2.3 ALE Approximation

ALE is estimated from the available dataset instances. (Apley and Zhu, 2020) proposed to divide the feature domain in K bins of equal size and to estimate the local effects in each bin by evaluating the model f at the bin limits:

$$\hat{f}^{ALE}(x_s) = \sum_{k=1}^{k_x} \frac{1}{|\mathcal{S}_k|} \sum_{i: \mathbf{x}^i \in \mathcal{S}_k} \left[ f(z_k, \mathbf{x}_c^i) - f(z_{k-1}, \mathbf{x}_c^i)) \right]$$
(2)

where  $k_x$  the index of the bin that  $x_s$  belongs to, i.e.  $k_x: z_{k_x-1} \leq x_s < z_{k_x}$  and  $\mathcal{S}_k$  is the set of training instances of the k-th bin, i.e.  $\mathcal{S}_k = \{\mathbf{x}^i: z_{k-1} \leq x_s^i < z_k\}$ . (Gkolemis et al.) proposed the Differential ALE (DALE) that computes the local effects on the training instances using auto-differentiation:

$$\hat{f}^{\text{DALE}}(x_s) = \Delta x \sum_{k=1}^{k_x} \frac{1}{|\mathcal{S}_k|} \sum_{i: \mathbf{x}^i \in \mathcal{S}_k} f^s(\mathbf{x}^i)$$
 (3)

Their method has significant computational advantages and allows the recomputation of the accumulated effect with different bin-splitting with near-zero additional computational cost. Moreover, by avoiding the use of artificial samples at the bin limits DALE allows the use of wider bins without out-of-distribution sampling. [Maybe connect with our work]. In both cases, the approximations partition the feature domain in K equally-sized bins, without considering the underlying local effects.

# 3 Uncertainty-Aware ALE (UALE)

We define UALE and provide an interval-based formulation that partitions the feature domain into non-overlapping intervals (bin-splitting) and defines the bin-effect and the bin-uncertainty. Then, we present UALE approximation, formally proving the bin-splitting requirements for an unbiased estimation of the uncertainty. Based on this, we present an optimization method to determine an appropriate variable-width partitioning.

# 3.1 UALE Definition And Interval-Based Formulation

We quantify the uncertainty of the local effects at  $x_s=z$  with the standard deviation of the local explanations,  $\sigma(z)$ , where:

$$\sigma^{2}(z) = \mathbb{E}_{X_{c}|X_{s}=z} \left[ \left( f^{s}(z, X_{c}) - \mu(z) \right)^{2} \right]$$
 (4)

The uncertainty emerges from (a) the feature correlations and (b) the implicit feature interactions of the black-box function. We also define the accumulated uncertainty at  $x_s$ ,

Figure 1: UALE-concept-figure

as the accumulation of the standard deviation of the local effects:

$$f_{\sigma}^{\text{ALE}}(x_s) = \int_{x_s \min}^{x_s} \sigma(z) \partial z$$
 (5)

UALE defines the effect as a pair that consists of the average effect and the uncertainty  $(\mu(z), \sigma(z))$ . For visualizing it, we use the accumulated average effect, as defined in Eq. (1), and the accumulated uncertainty, as defined in Eq. (5). Figure 1 illustrates an example [Describe after putting the figure].

For the interval-based formulation, we define the bin-effect  $\mu(z_1, z_2)$  and the bin-uncertainty  $\sigma(z_1, z_2)$  as:

$$\mu(z_1, z_2) = \mathbb{E}_{z \sim \mathcal{U}(z_1, z_2)}[\mu(z)] = \frac{\int_{z_1}^{z_2} \mu(z) \partial z}{z_2 - z_1}$$
 (6)

$$\sigma^{2}(z_{1}, z_{2}) = \mathbb{E}_{z \sim \mathcal{U}(z_{1}, z_{2})}[\sigma^{2}(z)] = \frac{\int_{z_{1}}^{z_{2}} \sigma^{2}(z) \partial z}{z_{2} - z_{1}}$$
 (7)

Intuitively, if we randomly draw a point  $z^*$  from a uniform distribution  $z^* \sim \mathcal{U}(z_1, z_2)$ , the bin-effect (Eq. (6)) and the bin-uncertainty (Eq. (7)) are the expected average effect and the expected uncertainty, respectively. Also, if we denote as  $\mathcal{Z}$  the sequence of K+1 points that partition the domain of the s-th feature into K variable-size intervals, i.e.  $\mathcal{Z} = \{z_0, \ldots, z_K\}$ , we can reformulate UALE to:

$$\tilde{f}_{\mathcal{Z},\mu}^{\text{ALE}}(x_s) = \sum_{k=1}^{k_x} \mu(z_{k-1}, z_k)(z_k - z_{k-1})$$
 (8)

$$\tilde{f}_{Z,\sigma}^{ALE}(x_s) = \sum_{k=1}^{k_x} \sigma(z_{k-1}, z_k)(z_k - z_{k-1})$$
 (9)

where  $k_x$  is the index of the bin such that  $z_{k_x-1} \le x_s < z_{k_x}$ . Eq. (8) and Eq. (9) are the interval-based formulation of UALE.

### 3.2 UALE Interval-Based Approximation

For approximating the bin-effect and the bin-uncertainty, we use the set  $\mathcal{S}_k$  of dataset instances with the s-th feature lying inside the k-th bin, i.e.,  $\mathcal{S}_k = \{\mathbf{x}^i : z_{k-1} \leq x_s^i < z_k\}$ . The bin-effect is approximated with,

$$\hat{\mu}(z_{k-1}, z_k) = \frac{1}{|\mathcal{S}_k|} \sum_{i: \mathbf{x}^i \in \mathcal{S}_k} \left[ f^s(\mathbf{x}^i) \right]$$
 (10)

which, as shown by(Gkolemis et al.), is an unbiased estimator of Eq. (6) under the assumption that the points are uniformly distributed inside the intervals. The approximation for the uncertainty (Eq. (7)) from the available instances is defined as follows:

$$\hat{\sigma}^{2}(z_{k-1}, z_{k}) = \frac{1}{|\mathcal{S}_{k}|} \sum_{i: \mathbf{x}^{i} \in \mathcal{S}_{k}} \left( f^{s}(\mathbf{x}^{i}) - \hat{\mu}(z_{1}, z_{2}) \right)^{2}$$
(11)

In the Appendix B, we show that  $\hat{\sigma}^2(z_1,z_2)$  is an unbiased estimator of  $\sigma^2_*(z_1,z_2) = \frac{\int_{z_1}^{z_2} \mathbb{E}_{\mathbf{x}_c|x_s=z} \left[ (f^s(z,\mathbf{x}_c) - \mu(z_1,z_2))^2 \right] \partial z}{z_2-z_1}$ . In Theorem 3.1 we prove that in the general case,  $\sigma^2_*(z_1,z_2) \geq \sigma^2(z_1,z_2)$ , so  $\hat{\sigma}^2(z_1,z_2)$  is an overestimation of the bin uncertainty.

**Theorem 3.1.** If we define (a) the residual  $\rho(z)$  as the difference between the expected effect at z and the bin-effect, i.e  $\rho(z) = \mu(z) - \mu(z_1, z_2)$ , and (b)  $\mathcal{E}(z_1, z_2)$  as the mean squared residual of the bin, i.e.  $\mathcal{E}(z_1, z_2) = \frac{\int_{z_1}^{z_2} \rho^2(z) \partial z}{z_2 - z_1}$ , then it holds

$$\sigma_*^2(z_1, z_2) = \sigma^2(z_1, z_2) + \mathcal{E}^2(z_1, z_2)$$
 (12)

*Proof.* The proof is at Appendix C.  $\Box$ 

We refer to  $\mathcal{E}^2(z_1, z_2)$  as bin-error. Based on Theorem3.1, the estimation is unbiased only when  $\mathcal{E}^2(z_1, z_2) = 0$ .

#### 3.3 Bin-Splitting

The quality of UALE approximation is affected by (a) the population of samples in each bin and (b) the error term  $\mathcal{E}(z_1,z_2)$  of each bin. On the one hand, we favor a partitioning of the feature domain in large-bins to allow a robust estimation of  $\hat{\mu}(z_1,z_2),\hat{\sigma}(z_1,z_2)$  from a sufficient population of samples. On the other hand, we want to minimize the cummulative bin-error, i.e.,  $\mathcal{E}_{\mathcal{Z}}^2 = \sum_{k=1}^K \mathcal{E}^2(z_1,z_2)\Delta z_k$ , where  $\mathcal{Z}=\{z_0,\cdots,z_K\}$  and  $\Delta z_k=z_k-z_{k-1}$ . For convenience, we refer to  $\mathcal{E}_{\mathcal{Z}}^2$  as accumulated error. We search for a partitioning  $\mathcal{Z}_*$  that balances this trade-off.

Due to the Corollary 3.1.1, minimizing  $\mathcal{E}_{\mathcal{Z}}^2$  is equivalent to minimizing  $\sigma_*^2(z_{k-1}, z_k)\Delta z_k$ , which can be directly estimated by  $\hat{\sigma}^2(z_{k-1}, z_k)\Delta z_k$ .

**Corollary 3.1.1.** A bin-splitting Z minimizes the accumulated if and only if it also minimizes the accumulated error.

$$\underset{\mathcal{Z} = \{z_0, \dots, z_K\}}{\arg \min} \mathcal{E}_{\mathcal{Z}}^2 = \underset{\mathcal{Z} = \{z_0, \dots, z_K\}}{\arg \min} \sum_{k=1}^K \sigma_*^2(z_1, z_2) \Delta z_k \quad (13)$$

*Proof.* The proof is at Appendix D

Based on the above, we set-up the following optimization problem:

$$\min_{\mathcal{Z}=\{z_0,\dots,z_K\}} \quad \mathcal{L} = \sum_{k=1}^K \tau_k \hat{\sigma}^2(z_{k-1}, z_k) \Delta z_k$$
where 
$$\Delta z_k = z_k - z_{k-1}$$

$$\tau_k = 1 - \alpha \frac{|S_k|}{N}$$
s.t. 
$$|S_k| \ge N_{\text{NPB}}$$

$$z_0 = x_{s,min}$$

$$z_K = x_{s,max}$$
(14)

The objective  $\mathcal{L}$  searches for a partitioning  $\mathcal{Z}_* = \{z_0,\dots,z_K\}$  with low aggregated error  $\mathcal{E}^2_{\mathcal{Z}}$ . In case of many partitionings with similar aggregate-error, the term  $\tau_K$  favors the partitioning with more points per bin (with wider bins). The constraint of at least  $N_{\text{PPB}}$  points per bin sets the lowest-limit for a *robust* estimation. The user can choose to what extent they favor the creation of wide bins through (a) the parameter  $\alpha$  that controls the discount  $\tau_k$  and (b) the parameter  $N_{\text{PPB}}$  that sets the minimum population per bin. For providing a rough idea, in our experiments we set  $\alpha=0.2$  which means that the discount ranges between [0%,20%] and  $N_{\text{PPB}}=\frac{N}{20}$ , where N is the dataset size.

For solving the optimization problem of Eq.14 we add some constraints. First, we set a threshold  $K_{max}$  on the maximum number of bins which, in turn, defines the minimum bin-width, i.e.  $\Delta x_{min} = \frac{x_{s,max} - x_{s,min}}{K_{max}}$ . Based on that, we restrict the bin-limits to be multiples of the minimum width, i.e.  $z_k = k \cdot \Delta x_{min}$ , where  $k \in \{0, \cdots, K_{max}\}$ . In this discretized solution space, we find the global optimal using dynamic programming.

### 3.3.1 Solve Bin-Splitting with Dynamic Programming

# 4 SIMULATION EXAMPLES

In XAI it is common (Aas et al., 2021; Herbinger et al., 2022) to use simulation examples for compering methods, because the data-generating distribution  $p(\mathbf{X})$  and the predictive function  $f(\cdot)$  are known, and, therefore, the ground-truth is accesible. We follow this common XAI practice and we split the evaluation in two groups of examples. The first group (Section 4.1) compares UALE definition against PDP-ICE in modeling the average effect and the uncertainty. The second group (Section 4.2) compares UALE approximation using the automatic bin-splitting against the fixed-size alternative, in estimating the average effect and the uncertainty.

#### 4.1 Case 1: UALE vs PDP-ICE

In this simulation, we compare UALE against PDP-ICE. The aim of this example is to highlight that when features are correlated, PDP and ICE fail in quantifying the average effect and the heterogeneity, even when the black-box function is relatatively simple.

We use the following distribution:  $p(\mathbf{x}) = p(x_3)p(x_2|x_1)p(x_1)$  where  $x_1 \sim \mathcal{U}(0,1)$ ,  $x_2 = x_1$  and  $x_3 \sim \mathcal{N}(0,\sigma_3^2)$  and the following piecewise linear function:

$$f(\mathbf{x}) = \begin{cases} f_{\text{lin}} + \alpha f_{\text{int}} & \text{if } f_{\text{lin}} < 0.5\\ 0.5 - f_{\text{lin}} + \alpha f_{\text{int}} & \text{if } 0.5 \le f_{\text{lin}} < 1 \\ \alpha f_{\text{int}} & \text{otherwise} \end{cases}$$
(15)

where  $f_{1\text{in}}(\mathbf{x}) = a_1x_1 + a_2x_2$  and  $f_{\text{int}}(\mathbf{x}) = x_1x_3$ . The linear term  $f_{1\text{in}}$  includes the two correlated features and the interaction term  $f_{\text{int}}$  the two non-correlated. We evaluate the effect computed by UALE and PDP-ICE in three cases; (a) without interaction ( $\alpha=0$ ) and equal weights ( $a_1=a_2$ ), (b) without interaction ( $\alpha=0$ ) and different weights ( $a_1\neq a_2$ ) and (c) with interaction ( $\alpha\neq 0$ ) with equal weights ( $a_1=a_2$ ). In all cases, we provide the ground-truth average effect and uncertainty computed with analytical derivations (proofs at the Appendix) and we compare it against the approximations provided by each method, after generating N=500 samples.

No Interaction, Equal weights. In this set-up,  $\alpha=0$  (no interaction) and the  $a_1=a_2=1$  (equal weights). The ground truth effect  $f_{\mu}^{\rm GT}(x_1)$  is:  $x_1$  in [0,0.25],  $-x_1$  in [0.25,0.5] and 0 in [0.5,1]. The uncertainty is  $f_{\sigma^2}^{\rm GT}(x_1)=0$  in [0,1] because  $x_1$  does not interact with ohter variables. In Figure 2, we observe that PDP effect is wrong and ICE plots imply the existence of heterogeneous effects. In contrast, UALE models correctly both the average effect and the zero uncertainty and approximates it accurately finding a zero aggregate residual error partitioning of  $x_1$  domain.

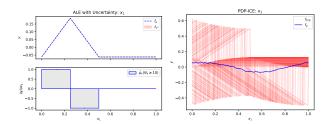


Figure 2: No interaction, Equal weights: Feature effect for  $x_1$  using UALE (Left) and PDP-ICE (Right).

**No Interaction, Non-Equal Weights.** In this set-up,  $\alpha = 0$  (no interaction),  $a_1 = 2$  and  $a_2 = 0.5$  (non-equal

weights). The non-equal weights have implications at both the gradient and the interval of the piece-wise linear regions, i.e.,  $f_{\mu}^{\rm GT}(x_1)$  is:  $2x_1$  in [0,0.2],  $-2x_1$  in [0.2,0.4] and 0 in [0.4,1]. As before, the ground-truth uncertainty is  $f_{\sigma^2}^{\mathcal{GT}}(x_1)=0$  because  $x_1$  does not interact with other features. In Figure 3, we observe that PDP estimation is completely opposite to the ground-truth effect, i.e. negative in the region [0,0.2), positive in [0.2,0.4), and the ICE erroneously implies the existence of heterogeneous effects. As before, ALE quantifies correctly the ground truth effect, the zero-uncertainty and partioning correctly the  $x_1$  domain.

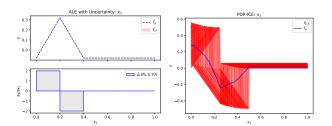


Figure 3: No interaction, Different weights: Feature effect for  $x_1$  using UALE (Left) and PDP-ICE (Right).

Uncertainty, Equal weights. In this set-up, we activate the interaction term, i.e. a=1, keeping the weights equal  $a_1=a_2=1$ . The interaction term provokes heterogeneous effects for features  $x_1$  and  $x_3$ , because the local effects at  $x_1$  depend on the unknown value of  $X_3$  and vice-versa. The effect of  $x_2$  has zero-uncertainty since it does not appear in any interaction term. As we observe in Figure 4, UALE correctly models the average effect and the uncertainty in all cases. On the other hand, PDP-ICE quantifies correctly the average effect and the uncertainty only in the case of feature  $x_3$ , because  $x_3$  is independent from other features. For the correlated features ( $x_1, x_2$ ) both the average effect (PDP) and the uncertainty (ICE) are wrong.

**Discussion.** Despite the model's simplicity, PDP-ICE fail in modeling both the average effect and the uncertainty. The error is due to the method's formulation, i.e. the use of the marginal distribution that ignores correlations between features, and not due to poor approximation because of limited samples. In contrast, UALE models both correctly and estimates them accurately extracting the regions with constant effect. The examples above do not cover the case of an interaction term between correlated features, for example a term  $x_1x_2$ , because there is an open debate about the ground-truth effect in this case (Grömping, 2020).

#### 4.2 Case 2: Bin-Splitting

In this simulation, we illustrate the advantages of automatic bin-splitting against the fixed-size alternative. For this reason, we generate a very big dataset applying dense sampling  $(N=10^6)$  and we treat the estimation with dense

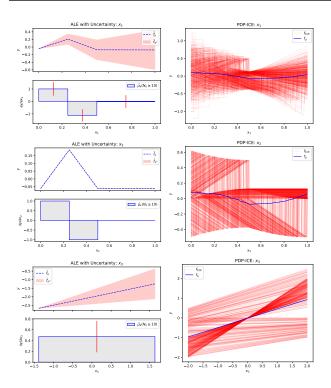


Figure 4: With interaction, equal weights: From top to bottom, feature effect for features  $\{x_1, x_2, x_3\}$  using UALE (left column) and PDP-ICE (right column).

fixed-size bins  $(K=10^3)$  as the ground-truth UALE. Afterwards, we generate less samples (N=500) and we compare the fixed-size estimation (for many different K) against UALE automatic bin-splitting. In all set-ups, we sample from  $p(\mathbf{x}) = p(x_2|x_1)p(x_1)$  where  $x_1 \sim \mathcal{U}(0,1)$  and  $x_2 \sim \mathcal{N}(x_1,\sigma_2^2=0.5)$ . We denote as  $\mathcal{Z}^{\mathrm{DP}}=\{z_{k-1}^{\mathrm{DP}},\cdots,z_K^{\mathrm{DP}}\}$  the sequence obtained by automatic bin-splitting and with  $\mathcal{Z}^{\mathrm{K}}$  the fixed-size splitting with K bins. The evaluation metrics we report are the average number of t=30 independent runs, using each time N=500 different samples.

**Metrics** The evaluation is done with regard to two metrics counting the mean error per bin, where the bin error is definde as the absolute difference of the approximation from the ground-truth. The first metric, Eq. (16), counts the mean bin error wrt average effect and the second, Eq. (17) the mean bin error wrt uncertainty:

$$\mathcal{L}_{\text{DP}|K}^{\mu} = \frac{1}{|\mathcal{Z}^{\text{DP}|K}|} \sum_{k \in \mathcal{Z}^{\text{DP}|K}} |\mu(z_k - z_{k-1}) - \hat{\mu}(z_k - z_{k-1})|$$
(16)

$$\mathcal{L}_{\text{DP}|K}^{\sigma} = \frac{1}{|\mathcal{Z}^{\text{DP}|K}|} \sum_{k \in \mathcal{Z}^{\text{DP}|K}} |\sigma(z_k - z_{k-1}) - \hat{\sigma}(z_k - z_{k-1})|$$
(17)

The ground truth  $\mu_k$   $(\sigma_k)$  is the mean value of the bineffects (bin-uncertainties) of the bins inside the interval defined by the wide bin. For better interpretation of the uncertainty error, we also provide the mean (per bin) error term  $\mathcal{L}^{\rho}_{\mathrm{DP}|\mathrm{K}} = \frac{1}{|\mathcal{Z}^{\mathrm{DP}|\mathrm{K}}|} \sum_{k \in \mathcal{Z}^{\mathrm{DP}|\mathrm{K}}} \mathcal{E}(z_{k-1}, z_k)$ .

Piecewise-Linear Function. In this set-up,  $f(\mathbf{x}) = a_1x_1 + x_1x_2$  is a piecewise-linear function with 5 different-width regions, i.e.,  $a_1$  equals to  $\{2, -2, 5, -10, 0.5\}$  in the intervals defined by the sequence  $\{0, 0.2, 0.4, 0.45, 0.5, 1\}$ . As we observe, in Figure 5, UALE's approximation is better than all fixedsize alternatives, in both mean effect ( $\mathcal{L}^{\mu}$ ) and uncertainty  $(\mathcal{L}^{\sigma})$ . For understanding the importance of variable-size splitting, we analyze the fixed-bin error. In the top-right of Figure 5, we observe the Law of Large Numbers (LLN);  $\mathcal{L}^{\mu}_{K}$  and K have positive correlation because as  $\Delta x$  becomes smaller less points contributing to each estimation. The interpretation of  $\mathcal{L}^{\sigma}$  is more complex. Given that the smallest piecewise-linear region is  $\Delta x = 0.05$  any K that is not a multiple of  $\frac{1}{\Delta x}=20$ , adds nuisance uncertainty  $\mathcal{L}_{\rm K}^{\rho}>0$  leading to a biased uncertainty approximation. For these K, uncertainy error decreases for bigger K because the length of the bins that lie in the limit between two piecewise linear regions becomes smaller. For K that are multiple of 20, i.e.  $K = \{20, 40, 60, 80, 100\}$ , the uncertainty approximation is unbiased  $\mathcal{L}_{\mathtt{K}}^{\rho} \approx 0$ . Here, as in  $\mathcal{L}^{\mu}$ , uncertainty estimation is worse as K grows larger (LLN). As we observe in the top left of Figure 5, automatic bin-splitting separates the fine-grain bins, e.g. regions [0.4, 0.45], [0.45, 0.5], and unites constant-effect regions into a single bin, e.g. region [0.5, 1].

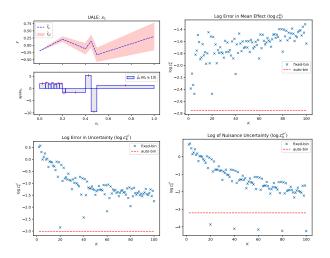


Figure 5: Figure 1

**Non-Linear Function.** In this set-up,  $f(\mathbf{x}) = 4x_1^2 + x_2^2 + x_1x_2$ , so the effect is non-linear in  $0 \le x_1 \le 1$ . In this set-up, wide bins increase  $\mathcal{L}^{\rho^2}$  making the uncertainty approximation more biased and narrow bins lead to a worse approximation. In Figure 6, we observe that automatic bin splitting manages to compromise the conflicting objectives, [Explain better]

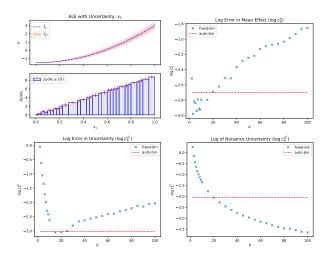


Figure 6: Figure 1

# 5 REAL-WORLD EXAMPLE

Here, we aim at demonstrating the usefuleness of uncertainty quantification and the advantages of automatic bin-splitting, on the real-world California Housing dataset (Pace and Barry, 1997).

ML setup The California Housing is a largely-studied dataset with approximately 20000 training instances, making it appropriate for robust Monte-Carlo approximations. The dataset contains D=8 numerical features with characteristics about the building blocks of California, e.g. latitude, longitude, population of the block or median age of houses in the block. The target variable is the median value of the houses inside the block in dollars that ranges between  $[15,500]\cdots 10^3$ , with a mean value of  $\mu_Y\approx 201\cdot 10^3$  and a standard deviation of  $\sigma_Y\approx 110\cdot 10^3$ .

We exclude instances with missing and outlier values. As outlier we define the feature values which ar over three standard deviations away from the mean feature value. We also normalize all features to zero-mean and unit standard deviation. We split the dataset into  $N_{tr}=15639$  training and  $N_{test}=3910$  test examples (80/20 split) and we fit a Neural Network with 3 hidden layers of 256, 128 and 36 units respectively. After 15 epochs using the Adam optimizer with learning rate  $\eta=0.02$ , the model achieves a normalized mean squared error (R-Squared) of 0.25 (0.75), which corresponds to a MAE of  $37 \cdot 10^3$  dollars.

Below, we illustrate the feature effect for three features: latitude  $x_2$ , population  $x_6$  and median income  $x_8$ . The particular features cover the main FE cases, e.g. positive/negative trend and linear/non-linear curve, and they are appropriate for illustration purposes. The complete results for all features, along with in-depth information about the preprocessing, training and evaluation parts are provided in the Appendix.

**Uncertainty Quantification** In real-world datasets, it is infeasible to obtain the ground truth FE for seamlessly evaluating the competitive methods. We selected the particular experiment, because, in broad terms, UALE and PDP-ICE plots aggree in the estimation of the average effect and uncertainty. Therefore, we can focus on juding the quality of the information provided by the two methods.

In Figure 7, we observe, from top to bottom, the effects for the latitude, population and the median income. The effect of UALE and PDP-ICE are similar for the population and the median income. The population has a negative impact that progressively decreases: from 400 to 1500 people the house value decreases with a rate of  $-150(\pm 140)$  dollars per added person, a rate that decreases from  $-80(\pm 80)$  to  $-60(\pm 60)$  dollars per added person as we move from 1500 to 2800 people. The level of uncertainty indicates significant variance in absolute value of the rate, but in the grant majority of instances the rate is negative. With the same inspection, we observe that the median annual income has a positive impact on the value (all numbers are thousands of dollars):  $20 \pm 15$  per 10 of added median income for incomes in [8, 15],  $32 \pm 20$  per 10 added income in [15, 60]and  $40 \pm 15$  per 10 added income in [60, 70]. The uncertainty indicates that there are less heterogeneous effects about the median income compared to the number of people. In both cases, we can end-up to the same conclusion by inspecting the PDP-ICE plots. For the latitude, there is a small difference in the explanations for the region [32, 35], where UALE estimates a less negative slope with less uncertainty than PDP, while the explanations are similar for the range [35, 39], where both methods reveal an increase in the uncertainty around the feature value 37.5.

In general, we observe that UALE complements ALE in the quantification of the heterogeneous effects, similarly to as ICE complements PDP. The automatic extraction of constant-effect and constant-uncertainty regions provided by UALE is helpful for an easier interpretation. On the other hand, ICE plots sometimes are more descritive on locating the type of heterogeneity, whereas UALE quantifies only the level (not the type) of heterogeneity.

**Bin Splitting** In this part, we evaluate the robustness of the approximation using automatic bin-splitting. Following the evaluation framework of Section 4.2, we treat as ground-truth the effects computed on the full training-set

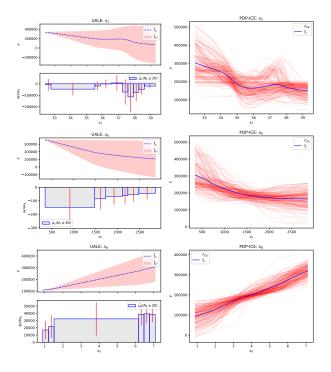


Figure 7: Figure 1

N=20000 with dense fixed-size bin-splitting (K=80). Given the big number of samples, we make the hypothesis that the approximation with dense binning is close to the ground truth. Afterwards, we randomly select less samples N=1000 and we compare UALE approximation with all possible fixed-size alternatives, repeating this process for 30 times for robust results. In Figure 8, we illustrate the mean values for  $\mathcal{L}^{\mu}$ ,  $\mathcal{L}^{\sigma}$  of the 30 repetitions. We observe that automatic bin-splitting provides (close to the) best approximation in the three features. In the Appendix, we provide the same evaluation for all features.

# 6 DISCUSSION

[Add Conclusion, limitations, and future work]

# Acknowledgments

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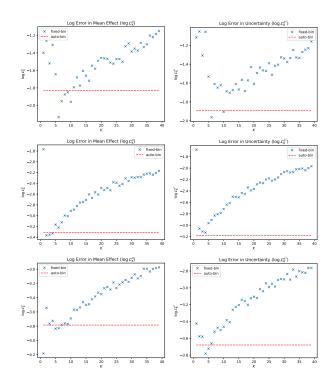


Figure 8: Figure 1

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# A List with Symbols

- **B** Show that is unbiased estimator
- C Proof of Theorem 3.1
- D Proof of Corollary 3.3
- E Dynamic Programming

[Review all the description] For achieving a computationally-grounded solution we set a threshold  $K_{max}$  on the maximum number of bins which also discretizes the solution space. The width of the bin can take discrete values that are multiple of the minimum step  $u = \frac{x_{s,max} - x_{s,min}}{K_{max}}$ . For defining the solution, we use two indexes. The index  $i \in \{0, \dots, K_{max}\}$  denotes the point  $(z_i)$  and the index  $j \in \{0, \dots, K_{max}\}$  denotes the position of the j-th multiple of the minimum step, i.e.,  $x_j = x_{s,min} + j \cdot u$ . The recursive cost function T(i,j) is

the cost of setting  $z_i = x_i$ :

$$\mathcal{T}(i,j) = \min_{l \in \{0,\dots,K_{max}\}} \left[ \mathcal{T}(i-1,l) + \mathcal{B}(x_l, x_j) \right]$$
(18)

where  $\mathcal{T}(0,j)$  equals zero if j=0 and  $\infty$  in any other case.  $\mathcal{B}(x_l,x_j)$  denotes the cost of creating a bin with limits  $[x_l,x_j)$ :

$$\mathcal{B}(x_l, x_j) = \begin{cases} \infty, & \text{if } x_j > x_l \text{ or } |\mathcal{S}_{(x_j, x_l)}| < N \\ 0, & \text{if } x_j = x_l \\ \hat{\sigma}^2(x_j, x_l), & \text{if } x_j \le x_l \end{cases}$$

$$(19)$$

The optimal solution is given by solving  $\mathcal{L} = \mathcal{T}(K_{max}, K_{max})$  and keeping track of the sequence of steps.

Discuss more aspects (e.g. Computational complexity)