

A Fast and Unified Partial Dependence Plot Algorithm

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Abstract

Partial Dependence Plots (PDPs) visualize how changes in a single feature affect the average model prediction. They are widely used in practice to interpret decision tree ensembles and other machine learning models. Joint-PDPs extend this idea to pairs of features, revealing their combined effect.

Partial Dependence Interaction Values (PDIVs) measure feature interactions. The Any-Order-PDIVs task computes these interactions for every feature subset across all rows of the dataset.

We introduce WOODELF++, a unified and efficient approach for computing all these useful explainability tools on decision tree ensembles, building on WOODELF, an algorithm for efficient SHAP computation. By deriving suitable metrics over pseudo-Boolean functions, WOODELF++ can compute PDPs (exact and approximate), Joint-PDPs, and Any-Order-PDIVs in a unified framework.

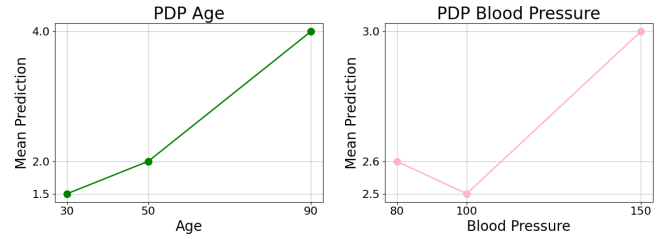
Our method delivers substantial complexity improvements over the state of the art, including an exponential gain for Any-Order-PDIVs. Additionally, we introduce and efficiently compute *Full PDPs*, which leverage the model’s split thresholds to faithfully capture its behavior across all possible feature values.

WOODELF++ is implemented in pure Python¹ and supports GPU acceleration. On a dataset with 400 000 rows, WOODELF++ computes PDP and Joint-PDP up to $6\times$ faster than the state of the art and up to five orders of magnitude faster than scikit-learn. For Any-Order-PDIVs, the gap is even larger: WOODELF++ computes all interaction values in 5 minutes, while the state of the art is estimated to require over 1 000 000 years.

1 Introduction

As decision tree ensembles are widely used in real-world applications [Blockeel *et al.*, 2023], understanding their predic-

¹WOODELF++ was migrated to the WOODELF GitHub package: <https://github.com/ron-wettenstein/woodelf>
Extended version of this work: <https://arxiv.org/abs/2605.14578>



(a) PDP on *age*, include three possible values: 30, 50 and 90. (b) PDP on *blood pressure*, with possible values: 80, 100 and 150.

Figure 1: PDP examples.

B[age] = 30 p1 = m.predict(B).mean()	B[bloodPressure] = 80 p1 = m.predict(B).mean()
B[age] = 50 p2 = m.predict(B).mean()	B[bloodPressure] = 100 p2 = m.predict(B).mean()
B[age] = 90 p3 = m.predict(B).mean()	B[bloodPressure] = 150 p3 = m.predict(B).mean()

Figure 2: Naive computation of the PDVs required to plot the PDPs in Fig. 1, carried out according to Def. 1.

tions is crucial for trust and decision-making [Yasodhara *et al.*, 2021; Ribeiro *et al.*, 2016]. This paper accelerates the computation of three widely used attribution metrics for decision tree ensembles: Partial Dependence Plot, Joint Partial Dependence Plot, and Partial Dependence Interaction Value.

1.1 Partial Dependence Plot (PDP)

A PDP [Friedman, 2001] illustrates how a model’s average prediction varies with changes in a specific feature. For example, Fig. 1a shows that the average prediction increases with *age*. PDPs are widely used in many scientific domains such as healthcare [Caruana *et al.*, 2015], ecology [Cutler *et al.*, 2007; Jibiki and Fukuda, 2024], and economics [Gnat, 2024].

A PDP is a line plot of several PDVs (Def. 1).

Definition 1 (Partial Dependence Value (PDV)). *A PDV measures the average prediction of a model M when a feature f_j is fixed to a given value v . For a dataset B :*

$$PDV_{v,f_j} = \frac{1}{|B|} \sum_{b \in B} M \left(\begin{cases} v & \text{if } f_i = f_j \\ b[f_i] & \text{otherwise} \end{cases} \right) \quad (1)$$

k : #selected values to plot; f : #features; F : Max number of features in a single tree; n : #samples (rows); T : #trees; D : tree depth; L : leaves per tree.

Figure 3: Notations.

To construct a PDP for feature f_j , we select k possible values of f_j , compute their PDVs, and plot them. The selected values are typically sampled from $B[f_j]$ or evenly spaced across its range.

For each feature f_a and samples $\{a_1, \dots, a_k\}$:

PDP requires $\forall_{1 \leq i \leq k} : PDV_{\{a_i\}, \{f_a\}}$

The naive way to construct these plots is to compute each PDV separately using Formula 1 (see Fig. 2). The overall complexity of computing PDPs for all features using this method is $O(kfnTD)$ (see Fig. 3 for notation). This naive approach remains the standard in widely used frameworks such as scikit-learn [Pedregosa *et al.*, 2011] and PiML [Sudjianto *et al.*, 2023].

[Friedman, 2001] also introduced a faster approach that approximates the PDVs, which we discuss in detail in Sec. 3.

Several extensions of PDPs have been proposed. Individual Conditional Expectation (ICE) [Goldstein *et al.*, 2015] plots can be viewed as PDPs computed for a single data row, revealing variation in feature effects across samples. Accumulated Local Effects (ALE) [Apley and Zhu, 2020] modifies the *PDV* computation by focusing on rows with feature values similar to v . A comparative analysis of PDPs, ALE, and Shapley values is presented in [Cook *et al.*, 2021]. Additionally, iPDP [Muschalik *et al.*, 2023] handles dynamic models that evolve over time, and StratPD [Parr and Wilson, 2020] takes a different approach: instead of interpreting the model, it generates plots that directly explain the training data.

1.2 Joint Partial Dependence Plot (Joint-PDP)

The concept of Partial Dependence Values extends naturally to feature sets: instead of fixing a single feature, we fix a set of features to chosen values - see Def. 2.

Definition 2 (Partial Dependence Value (PDV) on sets). *PDV on a subset of features $F = \{f_1, \dots, f_h\}$, set of values $V = \{v_1, \dots, v_h\}$ (write $V[f_i] = v_i$ for the value of f_i) and a dataset B is defined as follows:*

$$PDV_{V,F} = \frac{1}{|B|} \sum_{b \in B} M \left(\begin{cases} V[f_i] & \text{if } f_i \in F \\ b[f_i] & \text{otherwise} \end{cases} \right) \quad (2)$$

To visualize the joint contribution of a feature pair, it is useful to plot their PDVs. For features f_a, f_b with sampled values $\{a_1, \dots, a_h\}$ and $\{b_1, \dots, b_h\}$, Joint-PDP computes:

$$PDV_{\{a_i, b_j\}, \{f_a, f_b\}} \quad \text{for all } 1 \leq i, j \leq h$$

The naive approach, also used by scikit-learn, computes the *PDV* for every feature pair and value by fixing the pair to specific values and evaluating the model prediction (similar to Fig. 2, but fixing two features instead of one). This results in a time complexity of $O(k^2 f^2 nTD)$.

1.3 Partial Dependence Interaction Value (PDIV)

It is often useful to decompose the joint effect of features into their individual effects and the effect of their interactions. PDIVs (Def. 3) do exactly this.

Definition 3 (Partial Dependence Interaction Value (PDIV)). *A PDIV quantifies the interaction effect among the features $F = \{f_1, \dots, f_h\}$ when their values are fixed to $V = \{v_1, \dots, v_h\}$ (with $V[f_i] = v_i$). PDIVs on dataset B are defined as follows:*

$$PDIV_{V,F} = \sum_{S \subseteq F} (-1)^{|F|-|S|} \cdot PDV_{(V[f_i])_{\forall f_i \in S}, S} \quad (3)$$

For example: $PDIV_{\{4,1,6\}, \{f_2, f_5, f_8\}} = PDV_{\{4,1,6\}, \{f_2, f_5, f_8\}} - PDV_{\{4,1\}, \{f_2, f_5\}} - PDV_{\{1,6\}, \{f_5, f_8\}} - PDV_{\{4,6\}, \{f_2, f_8\}} + PDV_{\{4\}, \{f_2\}} + PDV_{\{1\}, \{f_5\}} + PDV_{\{6\}, \{f_8\}} - PDV_{\{\}, \{\}}$

Summing all interaction values over subsets of F reconstructs the PDV of F . Moreover, [Hiabu *et al.*, 2022] showed that the Shapley value of a feature f_i is a weighted sum of the PDIVs of all subsets S such that $f_i \in S$.

The *Any-Order-PDIVs* task computes PDIVs for each row c in the explained data C , across all feature subsets. Let $c|_S = \{c[f_i] : f_i \in S\}$, then the *Any-Order-PDIVs* task computes:

$$PDIV_{c|_S, S} \quad \text{for all } c \in C, S \subseteq F$$

An acceptable result for the *Any-Order-PDIVs* task is to return all non-zero PDIVs. A naive method runs predictions on the background data B (see Def 3, for simplicity, we assume $|B| = |C| = n$), for every feature subset $S \subseteq F$ (2^f subsets in total), and for every explained sample $c \in C$. This results in a total complexity of $O(2^f n^2 TD)$.

1.4 FastPD

FastPD [Liu *et al.*, 2025] is the current state-of-the-art tool for computing Any-Order-PDIVs. It exploits the structure and additivity of decision-tree ensembles to improve performance, achieving two main complexity improvements.

1. By processing each tree independently, FastPD reduces 2^f to 2^F . This is beneficial, as the number of features used in a tree (F) is typically much smaller than the total number of features (f).
2. FastPD preprocessing reduces the data-size term from $O(mn)$ to $O(m+n)$, where $m = |B|$ and $n = |C|$.

Overall, the complexity of FastPD for computing Any-Order-PDIVs is $O(T2^{D+F}(n+m))$. FastPD can also compute interaction values up to order s in complexity:

$$O \left(T \times n \times 2^D \times \sum_{i=0}^s \binom{F}{i} \right)$$

This complexity was not discussed in [Liu *et al.*, 2025]; we derived it ourselves and verified it empirically. Further details are provided in the supplementary material.

The GLEX package² implements the FastPD algorithm in R. Although FastPD theoretically supports separate background (B) and consumer (C) datasets, this option is not

²GLEX GitHub: <https://github.com/PlantedML/glex>

Task	scikit-learn	FastPD	WOODELF++
PDP, Sect. 3	$O(kfnTD)$	$O(T2^D F(n+k))$	$O(nTL + kTLD + TL2^D D)$
Any-Order-PDIVs, Sect. 5	-	$O(2^F nT)$	$O(n2^D TL + 4^D TL)$
Joint-PDP, Sect. 6	$O(k^2 f^2 nTD)$	$O(T2^D F^2(n+k^2 \log(f)))$	$O(k^2 f^2 \log(f) + nTL + k^2 \log(f) TLD^2 + TL2^D D^2)$

Table 1: Time complexity of SCIKIT-LEARN, FASTPD and WOODELF++ (see Fig. 3 for notations).

available in GLEX. For our experiments, we extended the package to support this setting while preserving its efficient implementation strategy. FastPD can also compute PDP in $O(T2^D F(n+k))$ and Joint-PDP in $O(T2^D F^2(n+k^2))$. This is done by computing first-order interactions for PDP and up to second-order interactions for Joint-PDP, and then reconstructing the PDVs from these interactions.

1.5 Contributions

Building on our previous work, WOODELF [Nadel and Wettenstein, 2026a], we formulate PDP, Joint-PDP, and Any-Order-PDIVs as cooperative-game metrics and introduce WOODELF++ for their efficient computation on decision tree ensembles. Our contributions include:

1. **PDP and Joint-PDP speedup:** We construct consumer datasets that enable efficient computation of PDP and Joint-PDP using local-attribution methods. We apply this approach to both WOODELF++ and FastPD, enhancing both methods. On large datasets, WOODELF++ is about $6\times$ faster than FastPD.
2. **Exponential complexity improvement for Any-Order-PDIVs:** In the worst case, where all splits use distinct features ($F = 2^D$), FastPD Any-Order-PDIVs complexity has a 2^{2^D} dependence, while WOODELF++ depends only on 4^D . We demonstrate this gap empirically: WOODELF++ completes Any-Order-PDIVs on 400 000 samples in just 5 minutes, whereas FastPD is estimated to take 1 000 000 years for this task.
3. **Full PDP:** We introduce a novel method for selecting the k plotted values in a PDP, yielding *full PDP* - a plot that displays the PDV for any possible feature value. Our speedups make it feasible to compute these plots for all model features within practical runtimes.
4. **PDP Approximation:** We demonstrate how WOODELF++ can be used to compute approximate PDPs, establishing a simple connection between approximate PDPs, exact PDPs, and Path-Dependence SHAP.

2 Preliminaries

2.1 Pseudo Boolean Functions

We remind the reader of Pseudo Boolean (PB) functions [Nadel, 2020] and their representation in Weighted Disjunctive Normal Form (WDNF) [Zhang and Jang, 2005], followed by the linearity property in this context.

Definition 4 (PB Function). A Pseudo Boolean (PB) function is a function of the form $F(x_1, \dots, x_h) : \{0, 1\}^h \rightarrow \mathbb{R}$.

Definition 5 (Positive and Negative Literal; Cube; WDNF). A literal is a Boolean variable x_i (positive literal) or its negation $\neg x_i$ (negative literal). A cube is a conjunction of literals. A Weighted Disjunctive Normal Form (WDNF) formula represents a PB function as:

$$F(x_1, \dots, x_h) = \sum_{k=1}^m w_k \cdot c_k(x_1, \dots, x_h),$$

where each c_k is a cube and $w_k \in \mathbb{R}$ is its weight.

Definition 6 (S_k, S_k^+, S_k^-). For a cube c_k , let S_k denote its set of variables, S_k^+ its positive literals, and S_k^- its negative literals.

Definition 7 (Linearity). A function g over WDNF formulas satisfies linearity if its value on a WDNF formula $F = \sum_{k=1}^m w_k \cdot c_k$ can be expressed as the weighted sum of its values on the individual cubes:

$$g(F) = \sum_{k=1}^m w_k \cdot g(c_k)$$

2.2 SHAP on Decision Trees and WOODELF

The inputs to Background SHAP are a decision tree ensemble M , a consumer dataset C containing the samples to be explained, and a background dataset B . For each consumer $c \in C$, the goal is to compute Shapley values [Shapley, 1953] for all of its features.

A cooperative game is defined over M , B , and c : when feature f_i participates ($f_i \in S$), it takes the consumer value $c[f_i]$; when it is missing ($f_i \notin S$), it is iteratively replaced with values from each baseline $b \in B$. The model prediction is evaluated for every $b \in B$, and the results are averaged.

WOODELF, introduced in our previous work [Nadel and Wettenstein, 2026a], models this game as a PB function F , assigning 1 to f_i when it participates and 0 when it is missing:

$$\frac{1}{|B|} \sum_{b \in B} M \left(\begin{cases} c[f_i] & \text{if } f_i \in S, \\ b[f_i] & \text{if } f_i \notin S \end{cases} \right) = F \left(\begin{cases} 1 & \text{if } f_i \in S, \\ 0 & \text{if } f_i \notin S \end{cases} \right) \quad (4)$$

WOODELF efficiently constructs a WDNF formula representing the PB function F . It then computes Background SHAP by applying a simple linear-time formula to the constructed WDNF. Path-Dependent SHAP is computed similarly, by constructing a slightly modified WDNF.

Importantly, WOODELF is not limited to Shapley values. This approach enables efficient computation of any function v that takes a WDNF formula as input and satisfies the linearity property (e.g., Banzhaf values).

A naive Background SHAP approach evaluates each consumer-background pair individually, resulting in $O(nm)$ complexity, where $|B| = m$, $|C| = n$, and L , T , and D are treated as constants. WOODELF and PLTreeSHAP [Zern *et al.*, 2023] use preprocessing to reduce this cost to $O(n + m)$.

More precisely, the complexity of WOODELF depends on the amortized cost of evaluating the function v across all cubes over the variables $\{x_1, \dots, x_D\}$ (where D is the tree depth), see Appendix D of [Nadel and Wettenstein, 2026b]. Let $O(v_c^{avg})$ denote the average computational cost of evaluating v on a single cube, and let $O(v_s)$ denote the number of unique feature subsets returned by v across all cubes. For example, in the case of Shapley values, each variable in a cube is considered individually, so $O(v_c^{avg}) = O(v_s) = O(D)$. The overall complexity of WOODELF is:

$$\text{Path-Dependent: } O(nTLv_s + TL3^D v_c^{avg})$$

$$\text{Background: } O(mTL + nTLv_s + TL3^D v_c^{avg})$$

3 Efficient PDP with WOODELF++

In this section, we extend WOODELF to efficiently compute all required PDVs for plotting the PDPs of all features.

The simple approach. Observe the strong correspondence between the PDV formula (Formula 1) and the model-prediction game formula (Formula 4): assigning $S = \{f_j\}$ and $c[f_j] = v$ in Formula 4 renders the expressions identical.

To better understand this correspondence, consider the PDV of *age* at $v = 50$. In the standard PDV computation, we set *age* to 50 for all samples in the background dataset B , run the model, and average the predictions (Fig. 2).

Equivalently, we can view this through a game-theoretic lens: imagine a consumer c with $c[\text{age}] = 50$ and the coalition where the feature *age* participates and is fixed to 50, while all other features are missing and filled with their values from B . The model’s expected prediction where only *age* participates is exactly the PDV of *age* at $v = 50$.

Leveraging this correspondence, WOODELF naturally extends to compute PDVs, yielding our WOODELF++. Let M be a decision tree ensemble, B its training set (or any dataset), and $\hat{f}$ the set of features in B . For each $f_i \in \hat{f}$, select k values $v_{1f_i}, \dots, v_{kf_i}$. We proceed as follows:

1. Construct a consumer dataset C with k rows. The t -th row of C contains the t -th selected value for each $f_i \in \hat{f}$.

$$C = ((v_{tf_i})_{\forall f_i \in \hat{f}})_{\forall 1 \leq t \leq k} \quad (5)$$

2. Use WOODELF with the background data B , the model M and the consumer data C , to produce a WDNF for each row in C .
3. For each WDNF formula F and feature f_i , apply:

$$PDV_{f_i} = F(f_i = 1, \forall f_{j \neq i} = 0) \quad (6)$$

For row t and feature f_i , this returns the average prediction when f_i participates (set to v_{tf_i}) and others are missing, i.e., $PDV_{v_{tf_i}, f_i}$.

Since all assignments (including $F(f_i = 1, \forall f_{j \neq i} = 0)$) are functions over WDNF formulas that satisfy linearity, WOODELF++ is guaranteed to compute them correctly.

The constructed consumer data C:

age	bloodPressure
30	80
50	100
90	150

The trainset is B.

The decision tree ensemble is M.

CPDVs = WOODELF(

M, C, B,

metric= For age: $F(\text{age}=1, \text{bloodPressure}=0) - F(\text{age}=0, \text{bloodPressure}=0)$

For bloodPressure: $F(\text{age}=0, \text{bloodPressure}=1) - F(\text{age}=0, \text{bloodPressure}=0)$

)

Figure 4: Using WOODELF++ to compute the CPDVs required for the plots in Fig. 1.

The efficient approach. Notice that Formula 6 above violates the *null player property*; that is, a variable whose truth value never affects the resulting weight might still receive a nonzero value. For example, the cube $3(\neg x_2)$ does not include the variable x_5 , yet it still affects PDV_{x_5} (since the assignment $\{x_5 = 1, \forall f_{j \neq 5} = 0\}$ satisfies it). As a result, a single cube may influence all the features in the dataset, which significantly hurts WOODELF’s performance.

Luckily, *Centered Partial Dependence Values* (CPDVs), obtained by subtracting the mean prediction on B from each PDV, do satisfy the null player property:

$$CPDV_{v, f_i} = PDV_{v, f_i} - \frac{\sum_{b \in B} M(b)}{|B|} \quad (7)$$

In the WDNF constructed by WOODELF, the mean term $\frac{1}{|B|} \sum_{b \in B} M(b)$ corresponds to the assignment where all variables are set to 0 (all features are “missing” and taken from B). Combining this insight with Formula 6 yields a simple formula for computing CPDV:

$$CPDV_{f_i} = F(f_i = 1, \forall f_{j \neq i} = 0) - F(\forall f_j = 0) \quad (8)$$

Fig. 4 illustrates how WOODELF++ and Formula 8 compute CPDVs efficiently.

Finally, Formula 9 is equivalent to Formula 8 on WDNFs. It satisfies both linearity and the null player property, making its use within WOODELF both correct and efficient:

$$CPDV_{f_i} = \sum_{k=1}^m w_k \times \begin{cases} 1 & \text{if } (S_k^+ = \{f_i\}) \wedge (f_i \notin S_k^-) \\ -1 & \text{if } (f_i \in S_k^-) \wedge (S_k^+ = \emptyset) \\ 0 & \text{otherwise} \end{cases} \quad (9)$$

The computation of CPDVs with the WOODELF package is straightforward — Fig. 5 shows the entire code. The computation is performed by calling WOODELF with a *CPDVMetric* instance, and requires no modifications to the package.

Recovering PDVs from CPDVs is immediate: add back the mean prediction on B . We now present the pseudo-code for our PDP algorithm. Alg. 1 receives a model M , a dataset B , and an integer k , and computes PDPs for all features in B , where each plot uses k sampled points.

A complexity analysis of Alg. 1 is provided in the supplementary material. In short, the WOODELF call (line 3) dominates the runtime: we show that $O(v_c^{avg}) = O(v_s) = O(D)$ and that cubes with $|S_k^+| > 1$ can be ignored, leaving only $O(2^D)$ relevant cubes. The complexity is:

$$\text{Exact PDP : } O(nTL + kTLD + TL2^D D)$$

```

1 class CPDVMetric(CubeMetric):
2     def calc_metric(self, s_plus: Set,
3         ↪ s_minus: Set):
4         if len(s_plus & s_minus) > 0:
5             return {}
6         pdp_values = {}
7         if len(s_plus) == 1:
8             for f in s_plus:
9                 pdp_values[f] = 1
10        if len(s_minus) == 0:
11            for f in s_minus:
12                pdp_values[f] = -1
13        return pdp_values
    
```

Figure 5: Python code for the class *CPDVMetric* that implements Formula 9 for a single cube with weight 1. It inherits WOODELF’s *CubeMetric*; a base class for metrics over WDNF formulas.

Algorithm 1 Efficient Partial Dependence Plot

- 1: **function** WPDP(M, B, k)
- 2: Select k possible values for each feature in B , and store them in C (see Formula 5).
- 3: $CPDVs = \text{WOODELF}(M, C, B, \text{CPDVMetric}())$
- 4: $PDVs = CPDVs + M(B).mean()$
- 5: Use $PDVs$ to plot the PDP of each feature.

Approximate PDP. Node cover is the number of training samples that reach each node during training. [Friedman, 2001] introduces a PDP approximation based on node cover. [Liu *et al.*, 2025] later observed that this uses the same approximation technique as Path-Dependent SHAP. In cooperative game terms, both methods are defined over the same game. [Liu *et al.*, 2025] also discusses an inconsistency of this approach: it may produce different explanations for functionally equivalent trees.

WOODELF++ leverages this connection to unify approximate and exact PDP under a single algorithm. Specifically, approximate PDP is obtained by computing the same PDV expression on the WDNF built with the Path-Dependent construction. In practice, this reduces to running the algorithm with an empty background dataset, so it defaults to the Path-Dependent setting (i.e., simply call: $\text{WPDP}(M, \emptyset, k)$).

The complexity of Alg. 1 with an empty background dataset (see the Path-Dependent complexity in Sect. 2.2):

$$\text{Approximate PDP} : O(kTLD + TL2^D D)$$

4 Full PDP with WOODELF++

Standard PDPs depend on both the number (k) and choice of sampled values. A small k can hide important patterns. For example, a TV-rating model may reveal a sharp viewer spike at lunch (11:30–13:00), but a PDP sampled every four hours (6:00, 10:00, 14:00, 18:00) would miss it.

Some effects are hard to capture even with large k . For instance, a fraud-detection model may learn that users who leave their annual salary at the default value of 60 000 are more likely to commit fraud. Even $k = 1000$ might miss this, as it occurs at a single value within a wide range.

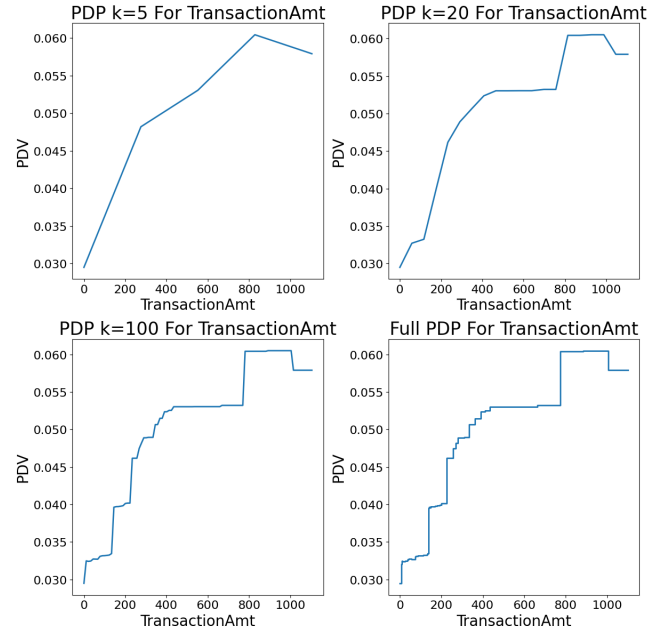


Figure 6: Comparison of PDPs for *TransactionAmt* on the IEEE-CIS dataset using $k = 5$, $k = 20$, and $k = 100$ uniformly spaced intervals versus the proposed *Full PDP*. Standard PDPs approximate the structure as k increases but fail to recover exact discontinuities even at $k = 100$. The Full PDP recovers the true step function using only 66 evaluation points corresponding to split thresholds.

To address this, we introduce *Full PDPs*, which capture all such behaviors in decision tree ensembles. For a given feature (e.g., *age*), we take all split thresholds from tree nodes that split on that feature and use them as x-axis values. The rationale is simple: the ensemble changes its prediction only when *age* crosses one of its thresholds. Between any two consecutive thresholds $th_1 < th_2$, the prediction remains constant:

$$\forall th_1 < v < th_2 (PDV_{v,age} = PDV_{th_1,age})$$

Thus, plotting only these points suffices.

Full PDPs reveal the complete piecewise-constant structure of the decision tree ensemble, including narrow spikes and single-value effects that sampled PDPs often miss. They are particularly useful for detecting overfitting and checking monotonicity. Fig. 6 compares sampled PDPs with different k values to a Full PDP, showing how the latter exactly recovers the true step function using only the feature’s split thresholds.

The main drawback is computational cost: large ensembles may have thousands of thresholds per feature. Previously, this made Full PDPs impractical. Both WOODELF++ (via Alg. 1) and FastPD handles large k efficiently, enabling Full PDPs at practical speeds.

5 Any-Order-PDIVs with WOODELF++

In this section, we extend WOODELF to compute all Partial Dependence Interaction Values (PDIVs; see Def. 3). Our approach follows three steps:

1. Derive a formula for PDIVs on a WDNF representation.

```

1 class PDIVMetric(CubeMetric):
2     INTERACTION_VALUE = True
3
4     def calc_metric(self, s_plus: Set,
5                     ↪ s_minus: Set):
6         if len(s_plus & s_minus) > 0:
7             return {}
8         pdivs = {}
9         for sm in all_subsets(s_minus):
10            s = tuple(s_plus | sm)
11            pdivs[s] = (-1) ** len(sm)
12        return pdivs
    
```

Figure 7: Python implementation of Formula 11 for a single cube with weight 1. `all_subsets` returns all subsets of a given set.

2. Use this formula to implement a Python class inheriting *CubeMetric* that computes all PDIVs of a cube.
3. Run WOODELF with that *CubeMetric*.

For a pseudo-Boolean function F and variable subset X , PDIV is defined by Formula 10 (combining Def. 2, Def. 3, and Formula 6):

$$PDIV_X = \sum_{S \subseteq X} (-1)^{|X|-|S|} F(\forall x_{i \in S} = 1, \forall x_{i \notin S} = 0) \quad (10)$$

Because PDIV is linear, computing $PDIV_X$ on the WDNF reduces to determining the contribution of each cube c_k separately and summing the results. We discard unsatisfiable cubes where $S_k^+ \cap S_k^- \neq \emptyset$. A cube affects only subsets X such that $S_k^+ \subseteq X \subseteq S_k$:

- If $X \not\subseteq S_k$, the interaction value is zero, since X contains at least one feature that does not appear in the cube and is therefore ignored. Any interaction involving an ignored feature is zero.
- If $S_k^+ \not\subseteq X$, at least one positive literal is falsified, so all assignments in Formula 10 return zero.

For subsets X where $S_k^+ \subseteq X \subseteq S_k$, Formula 10 sums the assignments of all $S \subseteq X$. The only such S to satisfy the cube and affect the summation is $S = S_k^+$. Thus:

$$PDIV_X = \sum_{k=1}^m w_k \times \begin{cases} (-1)^{|X|-|S_k^+|} & \text{if } S_k^+ \subseteq X \subseteq S_k \\ 0 & \text{otherwise} \end{cases} \quad (11)$$

The *PDIVMetric* code in Fig. 7 uses Formula 11 and computes the effect of a single cube on all relevant subsets X . Running WOODELF++ with *PDIVMetric* is straightforward:

```
return WOODELF(M, C, C, PDIVMetric())
```

This simple WOODELF++ call computes all the desired Any-Order-PDIVs. Here C is used as both consumer and background data, but a separate background dataset can be provided (e.g., using the training set as background). Limiting consumer size is recommended for memory efficiency. An empty background dataset produces approximate PDIVs via the Path-Dependent approach. As WOODELF is GPU-friendly, our method naturally supports GPU acceleration.

Complexity. For *PDIVMetric*, $O(v_s) = O(2^D)$ and $O(v_c^{avg}) = O((\frac{4}{3})^D)$, giving:

$$\text{Any-Order-PDIVs: } O(nTL2^D + TL4^D)$$

This complexity is exponentially better than the previous state of the art, and for balanced decision trees with 2^D nodes it becomes polynomial in the input size. A detailed complexity analysis is provided in the supplementary material.

The intuition behind this complexity improvement is as follows. The naive algorithm treats the model as a black box, leading to complexity exponential in the total number of features. FastPD [Liu *et al.*, 2025] improves on this by processing each tree separately; its complexity is exponential in the number of features used per tree, resulting in a worst-case dependence of 2^{2^D} , which remains exponential in the input size. WOODELF++ advances this line of improvement by treating each root-to-leaf path separately, yielding complexity exponential only in the number of features used per path, i.e., 2^D .

6 Joint-PDP with WOODELF++

To compute all Joint PDVs, we use the algorithm from Sect. 5 and the identity:

$$PDV_{\{a_i, b_j\}, \{f_a, f_b\}} = PDIV_{\{a_i, b_j\}, \{f_a, f_b\}} + PDIV_{\{a_i\}, \{f_a\}} + PDIV_{\{b_j\}, \{f_b\}} + PDIV_{\{\}, \{\}} \quad (12)$$

The algorithm:

1. Compute all PDIVs of order 2 and 1 using Sect. 5.
2. $PDIV_{\{\}, \{\}}$ is simply the average prediction on the background data. Compute it directly.
3. Apply Formula 12 to obtain PDVs for all feature pairs.

Deriving PDVs from PDIVs, as in Formula 12, is also used in FastPD [Liu *et al.*, 2025]. Since only up to second-order interactions matter, step 1 can ignore cubes with $|S_k^+| > 2$, leaving only $O(2^D D^2)$ cubes to consider.

To construct the consumer dataset, we sample k values per feature. For every pair of features f_a and f_b , the dataset must contain all k^2 value pairs (a_i, b_j) . A naive construction allocates k^2 rows per feature, resulting in $O(k^2 f)$ rows overall. Specifically, we first construct a size- k consumer data using the PDP procedure (Formula 5). Then, for every feature f_a and every sampled value a_i , we duplicate the constructed data C and set feature f_a to a_i in all rows:

$$\forall f_a \in F \forall 1 \leq i \leq k \forall c \in C : c[f_a] \leftarrow a_i$$

Using a more efficient construction (see supplementary material), we reduce the dataset size to only $O(k^2 \log(f))$ rows and adapt WOODELF++ to operate on this compressed representation. Overall complexity (see notations in Fig. 3):

$$O(k^2 f^2 \log(f) + nTL + k^2 \log(f) \cdot TLD^2 + TL2^D D^2)$$

This consumer data construction is not specific to WOODELF++. We also use it to accelerate FastPD on Joint-PDP.

As the PDIVs algorithm supports GPU acceleration, Joint-PDP is also GPU-friendly. Using full PDP sampling, one can plot *full Joint-PDPs* showing the joint dependence for all values of f_a and f_b .

7 Experimental Results

The PDP implementation is available in scikit-learn via the *partial_dependence* function. This function implements both the exact method (*method='brute'*) and the approximation method (*method='recursion'*) of PDV computation.

FastPD is implemented in the GLEX R package. It is the current state-of-the-art tool for computing Any-Order-PDIVs. The current implementation focuses on local feature attributions and supports only the case where the background and consumer datasets are identical. We extend the package to support distinct consumer and background datasets. Combined with our consumer-data constructions for PDP, Joint-PDP, and Full PDP, this enables efficient computation of these tasks within the FastPD framework.

Finally, we implemented WOODELF++ in Python for all discussed tasks and empirically validated it by comparing its outputs to those of scikit-learn.

We compare WOODELF++ running times against scikit-learn and FastPD on two large and widely used datasets:

- The IEEE-CIS fraud detection [kag, 2019] dataset with 472 432 rows and 397 features after one-hot encoding.
- The KDD Cup 1999 intrusion detection [kdd, 1999] dataset with 4 898 431 rows and 127 features after one-hot encoding.

For consistency, these are the same datasets we used in the original WOODELF paper [Nadel and Wettenstein, 2026a]. Additional experiments on smaller datasets and varying tree depths are reported in the supplementary material.

We used scikit-learn’s *HistGradientBoostingRegressor* model, which employs the histogram-based approach popularized by LightGBM [Ke *et al.*, 2017]. In all experiments, models were trained with 100 trees, each limited to a maximum depth of 6, matching XGBoost’s widely used default *max_depth*. Experiments were conducted on Google Colab. For CPU training, we enabled the high-memory runtime with 50GB RAM (the default is 12GB). For GPU training, we used the T4 runtime type with high-memory enabled. All experiment notebooks are available in the WOODELFEXPERIMENTS GitHub repository³.

As shown in Table 2, our approach outperforms the state of the art, with WOODELF++ achieving a $6\times$ speedup over FastPD on PDP and Joint-PDP. Notably, both FastPD and WOODELF++ significantly outperform scikit-learn, the most widely used tool for this task, with the gap becoming especially pronounced for large k : tasks that take hours or even several days using scikit-learn are completed by both approaches in under 10 minutes.

In Any-Order-PDIVs, where WOODELF++ offers an exponential complexity improvement over FastPD, the performance gap is substantial. On IEEE-CIS, this improvement translates to a reduction in runtime from an estimated one million years with FastPD to just five minutes using our approach.

³The WOODELFEXPERIMENTS repository containing all the notebooks used in our experiments: https://github.com/ron-wettenstein/WoodeLFExperiments/tree/main/woodeLF_experiments/woodeLF_IJCAI26

IEEE-CIS				
Task	scikit-learn	FastPD	WOODELF++	
			CPU	GPU
Exact PDP $k=5$	58.5 min	99 sec	14.5 sec	13.1 sec
Exact PDP $k=10$	112.4 min	99 sec	15.2 sec	10 sec
Exact PDP $k=100$	19 hours*	99 sec	14.8 sec	9 sec
Exact full PDP	59.3 min	98 sec	11.8 sec	7 sec
Exact Joint-PDP $k=5$	35 days*	130 sec	19 sec	25 sec
Approx. PDP $k=5$	3.3 sec	-	4.6 sec	5.4 sec
Approx. Joint-PDP $k=5$	19.4 min	-	10.2 sec	21.4 sec
Any-Order-PDIVs $n=10,000$	-	$6 \cdot 10^4$ years*	21 sec	22.6 sec
Any-Order-PDIVs	-	$2.9 \cdot 10^6$ years*	275 sec	30.7 sec

KDD Cup				
Task	scikit-learn	FastPD	WOODELF++	
			CPU	GPU
Exact PDP $k=5$	72.1 min	520 sec	95.6 sec	33.4 sec
Exact PDP $k=10$	100.2 min	519 sec	95.9 sec	33.5 sec
Exact PDP $k=100$	17 hours*	519 sec	96.2 sec	33.6 sec
Exact full PDP	116.5 min	519 sec	90.3 sec	26.8 sec
Exact Joint-PDP $k=5$	9 days*	568 sec	95.2 sec	37.5 sec
Approx. PDP $k=5$	5.5 sec	-	6.5 sec	8.7 sec
Approx. Joint-PDP $k=5$	10.1 min	-	21.6 sec	26 sec
Any-Order-PDIVs $n=10,000$	-	15 days*	80.6 sec	24 sec
Any-Order-PDIVs	-	20 years*	35 min	45 sec**

Table 2: Performance comparison between scikit-learn, FastPD, and WOODELF++. k is the number of sampled values per feature; n is the number of rows. Values marked with * are estimates; details on estimation are provided in the supplementary material. Dashes indicate unavailable or non-applicable measurements. **Due to RAM limitations, GPU evaluation of KDD “Any-Order-PDIVs” was conducted on an L4 runtime environment.

8 Conclusion

In this work, we extend our previous algorithm, WOODELF, originally designed for Shapley and Banzhaf values, into WOODELF++: a unified framework for efficiently computing PDPs, Joint-PDPs, and Any-Order PDIVs on decision tree ensembles. The two key ideas are to formulate these explanation tasks as linear metrics over WDNF formulas, and to construct compact consumer datasets that allow local-attribution algorithms to efficiently compute global dependence plots. *Full PDPs* take the latter idea a step further by constructing a consumer dataset that exactly capture the complete piecewise-constant behavior of decision tree ensembles.

This perspective yields substantial computational gains, especially on large datasets. For PDP and Joint-PDP, WOODELF++ is approximately $6\times$ faster than FastPD. For Any-Order-PDIVs, where WOODELF++ significantly improves the complexity, runtime decrease from several years to just a few minutes.

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