

The complexity of unsupervised learning of lexicographic preferences

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Unsupervised preferences learning

- User preferences are modeled by an order (total or not) on some set of items \mathcal{X}
- Preferences learning is useful in several practical applications, in particular when using recommendation
- Most preference learning approaches are supervised:
 - the learning data is a set of pairwise comparisons (A is preferred to B, etc.)
- This data can be costly to collect, in contrary to sales histories that are readily available
- Sales histories contain information about the preferences:
 - they mostly contain items that rank high in the preferences
- Unsupervised preferences learning uses a set of items chosen by a user as training set

Previous work and goal

Previous work by [FGM18]

- We proposed greedy unsupervised learning algorithms for lexicographic preferences models
- The approach is based on the idea that the user won't always chose her most preferred item (because this item is not available, because of a desire of variety, etc.)
- We rely on the following assumption:
the more preferred an item is, the more likely it is chosen,
i.e. if $o \succ o'$, then $p(o) \geq p(o')$
- The learning algorithms have some convergence properties and have been experimentally assessed, but their theoretical properties have not been studied

The goal of this new article is to study the time and sample complexity of unsupervised learning for several classes of lexicographic preferences models



Outline

- ① Context
- ② Lexicographic Preference Trees
- ③ Learning model

Toy example

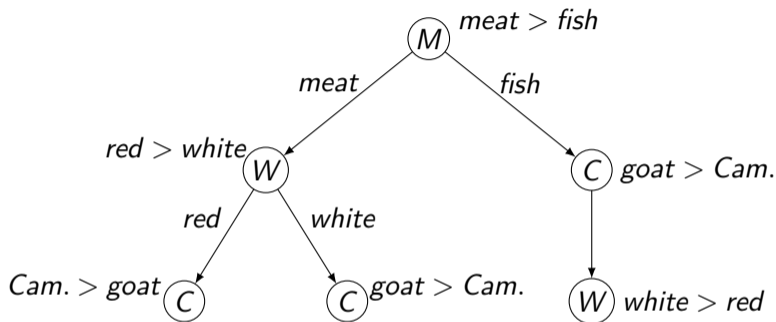
Imagine a user wants to book a dinner online. A menu is a tuple of values of three different attributes:

Main course (M) meat or fish

Wine (W) red or white

Cheese (C) goat or Camembert

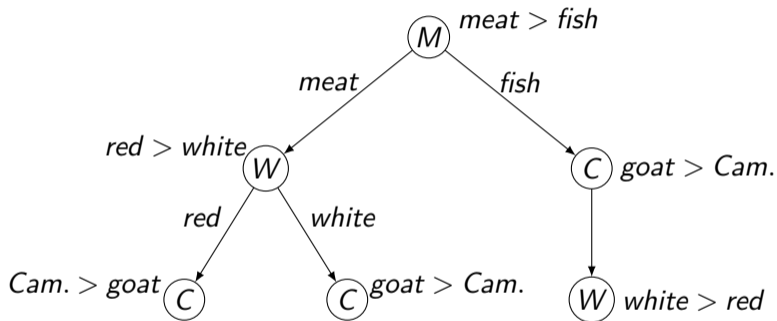
Lexicographic Preference Trees (LP-trees) [BCL⁺09]



LP-tree definition [BCL⁺09]

- Tree of attributes ordered by importance (root: most important)
- Edges can be labelled by a value or not
- Preferences rules associated to each node

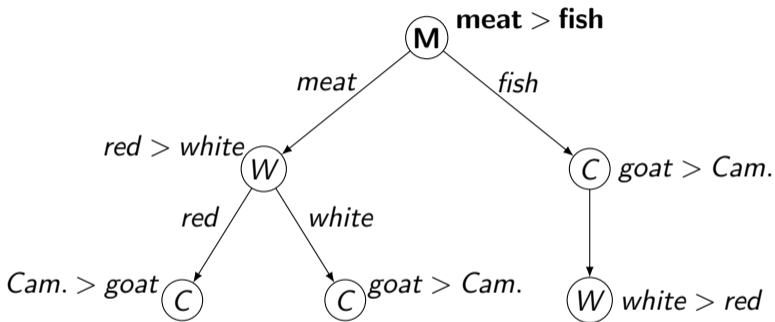
LP-trees semantics



We would like to compare:

meat – *red* – *goat* and *fish* – *white* – *goat*

LP-trees semantics

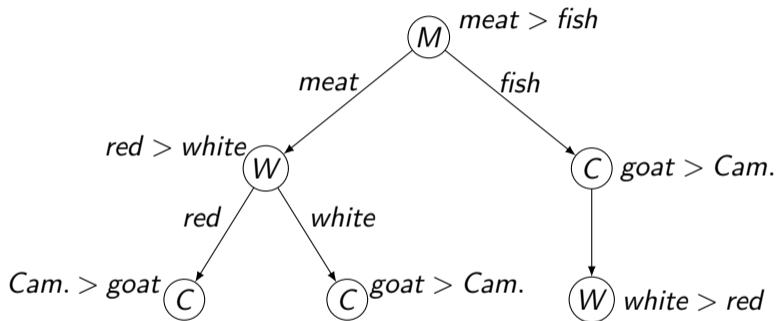


We would like to compare:

$meat - red - goat \succ fish - white - goat$

Any menu with meat is preferred to any menu with fish

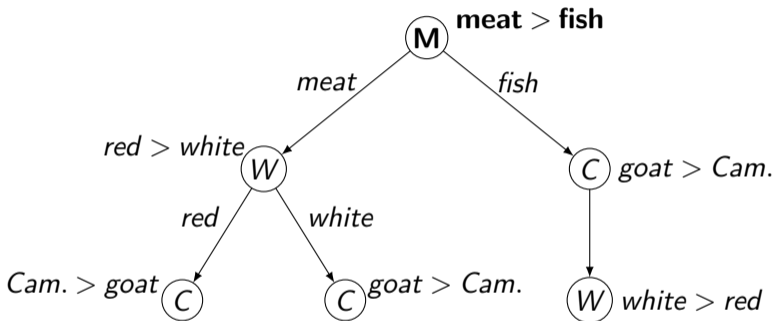
LP-trees semantics



We would like to compare:

meat – white – goat and *meat – white – Cam.*

LP-trees semantics

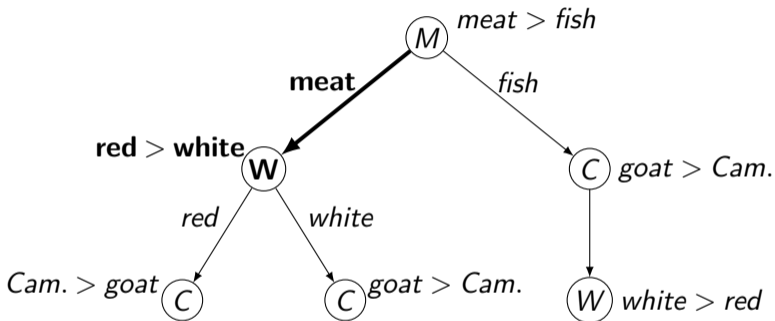


We would like to compare:

meat – white – goat and *meat – white – Cam.*

Root node can't decide the comparison

LP-trees semantics

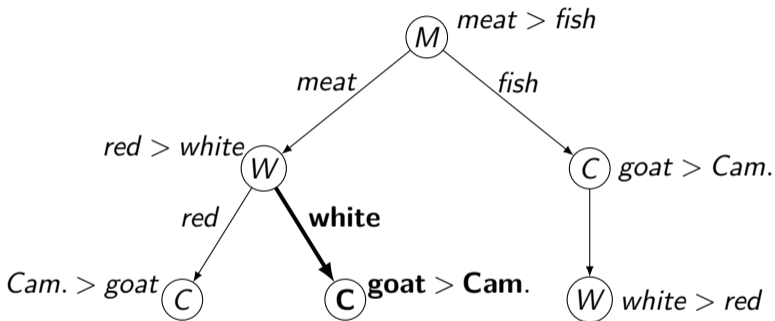


We would like to compare:

meat – white – goat and *meat – white – Cam.*

Among meat menus, any menu with red wine is preferred to any menu with white wine

LP-trees semantics



We would like to compare:

meat – white – goat \succ *meat – white – Cam.*



The LP-trees languages we study

Linear LP-trees with univariate nodes

One attribute per node, one branch only

Linear LP-trees with multivariate nodes

Multiple attributes per node, one branch only

LP-trees with multivariate nodes and a bounded number of leaves

Multiple attributes per node, max l branches

All these models represent total orders

Rank

- We consider a total relation \succ over $\underline{\mathcal{X}}$, a set of items
- The **rank** of an item $o \in \underline{\mathcal{X}}$, denoted as $rank(\succ, o)$, is defined as $1 +$ the number of elements preferred to o .
 - The rank of the preferred item is 1
 - The rank of the second preferred item is 2
 - ...
 - The rank of the least preferred item is $|\underline{\mathcal{X}}|$

Ranking loss

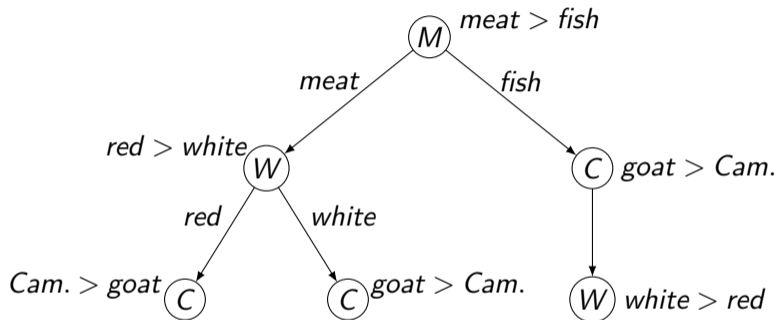
- To compare two models, we need a similarity measure but classic measures weight the same way the order on the preferred and the least preferred items
- In unsupervised learning, we use the ranking loss, the normalized difference of the mean rank of items drawn from p (preferred items are more weighted):

$$rloss_p(\gamma^*, \gamma') = \frac{1}{\underline{\mathcal{X}}} \left(\sum_{o \in \underline{\mathcal{X}}} p(o) rank(\gamma^*, o) - \sum_{o \in \underline{\mathcal{X}}} p(o) rank(\gamma', o) \right)$$

- Since γ^* (the ground truth) is unknown, the learning process seeks to minimize the mean rank of items drawn from the empirical distribution p_S according to γ' :

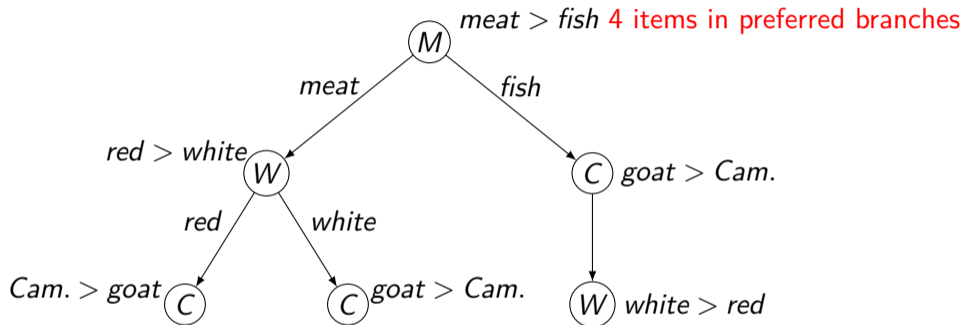
$$argmin_{\gamma'} \frac{1}{\underline{\mathcal{X}}} \sum_{o \in \underline{\mathcal{X}}} p_S(o) rank(\gamma', o)$$

The properties of the rank in a LP-tree



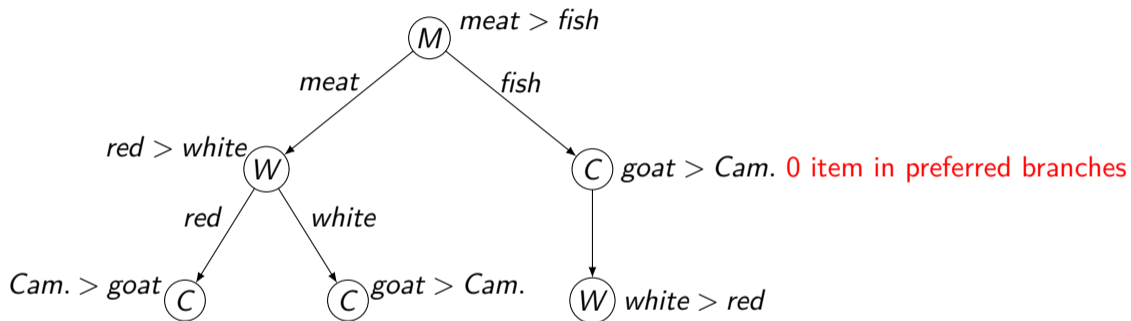
The rank of *fish/goat/red* is: 1 +

The properties of the rank in a LP-tree



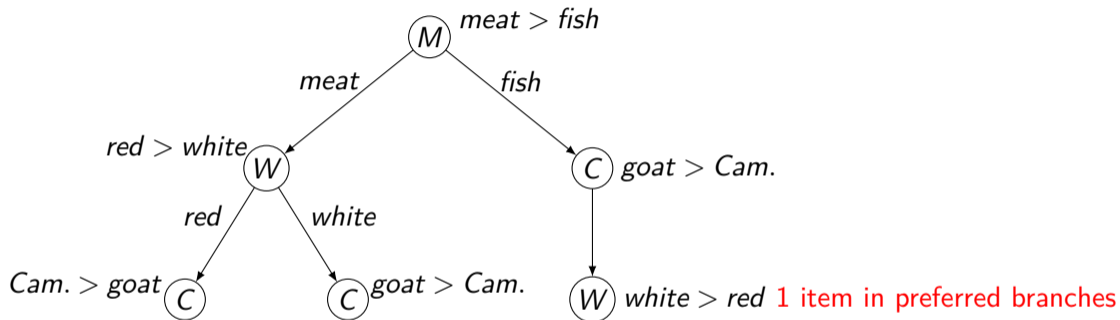
The rank of *fish/goat/red* is: 1 + 4 +

The properties of the rank in a LP-tree



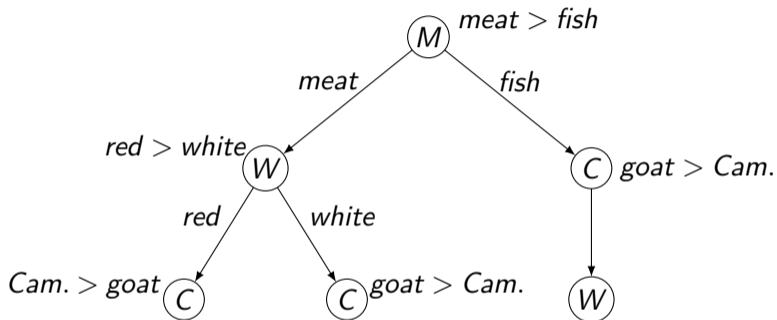
The rank of *fish/goat/red* is: $1 + 4 + 0 +$

The properties of the rank in a LP-tree



The rank of $fish/goat/red$ is: $1 + 4 + 0 + 1 = 6$

The properties of the rank in a LP-tree



The rank of *fish/goat/red* is: $1 + 4 + 0 + 1 = 6$

⇒ The rank of an element o can be decomposed into the sum of contributions of each node of one branch.

Rank decomposition

Formally:

$$\text{rank}(\phi, o) = 1 + \sum_{N \in \text{Nodes}(\phi, o)} |\underline{\text{Desc}}(N)| \times r(N, o[\text{Var}(N)])$$

where

- $\text{Nodes}(\phi, o)$ be the set of nodes in the branch compatible with o
- $|\underline{\text{Desc}}(N)|$ is the product of the sizes of the domain of the attributes below the node N
- $r(N, o[\text{Var}(N)])$ is the local rank (starting at 0) of the value of o for the attribute in N

The mean rank is:

$$\frac{1}{\underline{\mathcal{X}}} \sum_{o \in \underline{\mathcal{X}}} p_S(o) \text{rank}(\phi, o) =$$

$$\frac{1}{\underline{\mathcal{X}}} \left(1 + \sum_{N \in \text{Nodes}(\phi)} |\underline{\text{Desc}}(N)| \times p_S(\text{inst}(N)) \times E_{p_S}[r(N, o[\text{Var}(N)]) \mid \text{inst}(N)] \right)$$

Sample complexity definition

Let n be the number of attributes, $\hat{\phi}$ the unknown target LP-tree, ϕ^* the LP-tree with minimal mean rank (over a sample). We are looking for a function S such as, if there are at least $S(n, \delta, \epsilon)$ examples in the training set, then $Pr(rloss(\phi^*, \hat{\phi}) \leq \epsilon) \geq 1 - \delta$ (i.e., there is a high probability that the error is very small).

The result

Our upper bound on the sample complexity shows that it belongs in:

- $O(\frac{1}{\epsilon^2})$
- $O(\ln \frac{1}{\delta})$
- $O(\ln n)$
- $O(l^2)$
- $O(d^{4k})$,

where l = number of leaves, d = domain size,

k = max number of attributes per node, in practice $k < 4$

This complexity is low and shows that unsupervised learning is efficient

Sample Complexity : Sketch of Proof

$$\begin{aligned} \text{rloss}(\phi^*, \check{\phi}) &= \frac{1}{|\underline{\mathcal{X}}|} (E_p[\text{rank}(\phi^*, \cdot)] - E_p[\text{rank}(\check{\phi}, \cdot)]) \\ &\leq \frac{1}{|\underline{\mathcal{X}}|} (|E_p[\text{rank}(\phi^*, \cdot)] - E_{p_S}[\text{rank}(\phi^*, \cdot)]| \\ &\quad + E_{p_S}[\text{rank}(\phi^*, \cdot)] - E_{p_S}[\text{rank}(\check{\phi}, \cdot)] \\ &\quad + |E_{p_S}[\text{rank}(\check{\phi}, \cdot)] - E_p[\text{rank}(\check{\phi}, \cdot)]|) \\ &\leq \frac{2}{|\underline{\mathcal{X}}|} \max_{\phi \in \text{LPT}_I^k} |E_p[\text{rank}(\phi, \cdot)] - E_{p_S}[\text{rank}(\phi, \cdot)]| \end{aligned}$$

Sample Complexity : Sketch of Proof

$$\begin{aligned} & |E_p[\text{rank}(\phi, \cdot)] - E_{p_S}[\text{rank}(\phi, \cdot)]| \\ & \leq \left(\max_{\substack{N \in \text{nodes}(\phi) \\ v \in \underline{\text{Var}}(N)}} |p(v \wedge \text{inst}(N)) - p_S(v \wedge \text{inst}(N))| \right) \times \frac{d^k(d^k + 1)}{2} \times \sum_{N \in \text{nodes}(\phi)} \frac{|\text{Desc}(N)|}{|N|} \\ & \leq \left(\max_{v \subseteq \mathcal{X}, v \in \underline{\mathcal{V}}} |p(v) - p_S(v)| \right) \times \frac{d^k(d^k + 1)}{2} \times \sum_{N \in \text{nodes}(\phi)} \frac{|\text{Desc}(N)|}{|N|} \end{aligned}$$

where $d =$ bound on the domain size of the attributes.

Hoeffding's inequality is used to bound $|p(v) - p_S(v)|$

Linear lexicographic preference trees with univariate nodes

Order attributes in order of non-increasing

$$\text{Score}(p, X) = E_p^*(X) / (|\underline{X}| - 1)$$

\Rightarrow can be done in $O(n \log n |S| d \log d)$

Linear lexicographic preference trees with multivariate nodes

For every partition of \mathcal{X} :

- Order parts in order of non-increasing

$$\text{Score}(p, V) = E_p^*(V) / (|\underline{V}| - 1)$$

$\phi^* = \text{best partition}$

\Rightarrow exponential number of partitions



Time complexity for computing ϕ^*

Non-linear lexicographic preference trees \Rightarrow ???

Conclusion and future work

Conclusion

- Unsupervised learning relies on examples of chosen items and not on comparisons
- We focus on theoretical analysis of the time and sample complexity of unsupervised learning of three classes of lexicographic preferences trees

Future work

- We conjecture that the unsupervised learning of the optimal LP-tree is NP-hard
- We did not make any hypothesis on the distribution of p : we could study what law it follows in practice (maybe a power law distribution like the Zipf's law?)



Richard Booth, Yann Chevaleyre, Jérôme Lang, Jérôme Mengin, and Chattrakul Sombattheera.

Learning various classes of models of lexicographic orderings.

Preference learning, page 1, 2009.



Hélène Fargier, Pierre Francois Gimenez, and Jérôme Mengin.

Learning lexicographic preference trees from positive examples.

In *Proceedings AAAI 2018*, pages 2959–2966. ACM, 2018.