Online Learning of Weighted Relational Rules for Complex Event Recognition

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Abstract. Systems for symbolic complex event recognition detect occurrences of events in time using a set of event definitions in the form of logical rules. The Event Calculus is a temporal logic that has been used as a basis in event recognition applications, providing among others, connections to techniques for learning such rules from data. We advance the state-of-the-art by combining an existing online algorithm for learning crisp relational structure with an online method for weight learning in Markov Logic Networks (MLN). The result is an algorithm that learns complex event patterns in the form of Event Calculus theories in the MLN semantics. We evaluate our approach on a challenging real-world application for activity recognition and show that it outperforms both its crisp predecessor and competing online MLN learners in terms of predictive performance, at the price of a small increase in training time.

Keywords: Online Structure and Weight Learning · Markov Logic Networks · Event Calculus.

1 Introduction

Complex event recognition systems [7] process sequences of *simple events*, such as sensor data, and recognize *complex events* of interest, i.e. events that satisfy some temporal pattern. Systems for symbolic event recognition [3] typically use a knowledge base of first-order rules to represent complex event patterns. Learning such patterns from data is highly desirable, since their manual development is usually a difficult and error-prone task. The Event Calculus (EC) [23] is a temporal logical formalism that has been used as a basis in event recognition applications [2], providing among others, direct connections to machine learning, via Inductive Logic Programming (ILP) [8] and Statistical Relational Learning (SRL) [9].

Event recognition applications typically deal with noisy data streams [1]. Algorithms that learn from such streams are required to work in an online fashion, building a model with a single pass over the input [14]. Although a number of online relational learners have been proposed [12, 31, 18], learning theories in the EC is a challenging task that most relational learners, inluding the aforementioned ones, cannot fully undertake [28, 19]. As a result, two online algorithms have been proposed, both capable

of learning complex event patterns in the form of EC theories, from relational data streams. The first, OLED (Online Learning of Event Definitions) [21], adapts the Hoeffding bound-based [16] framework of [10] for online decision tree learning to an ILP setting. The second algorithm, OSLα (Online Structure Learning with Background Knowledge Axiomatization) [27], builds on the method of [18] for learning structure and weights in Markov Logic Networks (MLN) [29], towards online learning of EC theories in the MLN semantics.

Both these algorithms have shortcomings. OLED is a crisp learner, therefore its performance could be improved via SRL techniques that combine logic with probability. On the other hand, OSLα uses an efficient online weight learning technique, based on the AdaGrad algorithm [13], but its structure learning component is sub-optimal: It tends to generate large sets of rules, many of which are of low heuristic value with a marginal contribution to the quality of the learned model. The maintenance cost of such large rule sets during learning results in poor efficiency, with no clear gain in the predictive accuracy, while it also negatively affects model interpretability.

In this work we present a new algorithm that attempts to combine the best of these two learners: OLED's structure learning strategy, which is more conservative than OSLα's and typically explores much smaller rule sets, with OSLα's weight learning technique. We show that the resulting algorithm outperforms both its predecessors in terms of predictive performance, at the price of a tolerable increase in training time. We empirically validate our approach on a benchmark dataset of activity recognition.

The rest of this paper is structured as follows. In Section 2 we discuss related work, while in Section 3 we present some background material. In Section 3.1 we present OSLα and discuss its main limitations and in Section 4 we describe our proposed online structure and parameter learning method, after a brief presentation of the crisp version of OLED. In Section 5 we present our experimental evaluation, while in Section 6 we discuss some prospects of future work and conclude.

2 Related Work

Machine learning techniques for event recognition are attracting attention in the Complex Event Processing community [26]. However, existing approaches are relatively ad-hoc and they have several limitations [19], including limited support for background knowledge utilization and uncertainty handling. In contrast, we adopt an event recognition framework that allows access to well-established (statistical) relational learning techniques [8, 9], which can overcome such limitations. Moreover, using the EC allows for efficient event recognition via dedicated reasoners, such as RTEC [2].

However, learning with the EC is challenging for most ILP and SRL algorithms. A main reason for that is the non-monotonicity of the Negation as Failure operator that the EC uses for commonsense reasoning, which makes the divide-and-conquer-based search of most ILP algorithms inappropriate [28, 19, 20]. Non-monotonic ILP algorithms can handle the task [28, 5], but they scale poorly, as they learn whole theories from the entirety of the training data, while improvements to such algorithms that allow some form of incremental processing to enhance efficiency [4, 24] cannot handle data

arriving over time. Contrary to such approaches, OLED, the ILP algorithm we build upon, scales adequately and learns online [21].

A line of related work tries to "upgrade" non-monotonic ILP learners to an SRL setting, via weight learning techniques with probabilistic semantics [6, 11]. However, the resulting algorithms suffer from the same limitations, related to scalability, as their crisp predecessors. In the field of Markov Logic Networks (MLN), which is the SRL framework we adopt, OSLα is the sole algorithm capable of learning structure and weights for EC theories.

Online learning settings, as the one we assume in this work, are under-explored both in ILP and in SRL. A few ILP approaches have been proposed. In [12] the authors propose an online algorithm, which generates a rule from each misclassified example in a stream, aiming to construct a theory that accounts for all the examples in the stream. In [31] the authors propose an online learner based on Aleph[3] and Winnow [25]. The algorithm maintains a set of rules, corresponding to Winnow's features. Rules are weighted and are used for the classification of incoming examples, via the weighted majority of individual rule verdicts for each example, while their weights are updated via Winnow's mistake-driven weight update scheme. New rules are greedily generated by Aleph from missclassified examples. A similar approach is put forth by OSL [18], an online learner for MLN, which OSLα builds upon to allow for learning with the EC. OSL greedily generates new rules to account for misclassified examples that stream in, and then relies on weight learning to identify which of these rules are relevant. Common to these online learners is that they tend to generate unnecessarily large rule sets, which are hard to maintain. This is precisely the issue that we address in this work.

3 Background

We assume a first-order language, where predicates, terms, atoms, literals (possibly negated atoms), rules (clauses) and theories (collections of rules) are defined as in [9], while not denotes Negation as Failure. A rule is represented by $\alpha \leftarrow \delta_1, \ldots, \delta_n$, where α is an atom, (the head of the rule), and $\delta_1, \ldots, \delta_n$ is a conjunction of literals (the body of the rule). A term is ground if it contains no variables. We follow [9] and adopt a Prolog-style syntax. Therefore, predicates and ground terms in logical expressions start with a lower-case letter, while variable terms start with a capital letter.

The Event Calculus (EC) [23] is a temporal logic for reasoning about events and their effects. Its ontology consists of *time points* (integer numbers); *fluents*, i.e. properties that have different values in time; and events, i.e. occurrences in time that may alter fluents' values. The axioms of the EC incorporate the commonsense *law of inertia*, according to which fluents persist over time, unless they are affected by an event. We use a simplified version of the EC that has been shown to suffice for event recognition [2]. The basic predicates and its domain-independent axioms are presented in Table 1(a) and (b) respectively. Axiom (1) in Table 1(b) states that a fluent F holds at time T if it has been initiated at the previous time point, while Axiom (2) states that F continues to hold unless it is terminated. Definitions for initiatedAt/2 and terminatedAt/2 predicates are given in an application-specific manner by a set of *domain-specific* axioms.

[3] http://www.cs.ox.ac.uk/activities/machinelearning/Aleph/aleph

(a)

Predicate	**Meaning**
happensAt(E, T)	Event E occurs at time T.
initiatedAt(F, T)	At time T, a period of time for which fluent F holds is initiated.
terminatedAt(F, T)	At time T, a period of time for which fluent F holds is terminated.
holdsAt(F, T)	Fluent F holds at time T.

(b)

Domain-Independent Axioms

holdsAt$(F, T+1) \leftarrow$ (1)
  initiatedAt(F, T)

holdsAt$(F, T+1) \leftarrow$ (2)
  holdsAt(F, T),
  not terminatedAt(F, T)

(c)

Narrative for time 1:

happensAt$(walk(id_1), 1)$.
happensAt$(walk(id_2), 1)$.
holdsAt$(coords(id_1, 201, 454), 1)$.
holdsAt$(coords(id_2, 230, 440), 1)$
holdsAt$(direction(id_1, 270), 1)$
holdsAt$(direction(id_2, 270), 1)$

Annotation for time 1:

not holdsAt$(move(id_1, id_2), 1)$

Narrative for time 2:

happensAt$(walk(id_1), 2)$.
happensAt$(walk(id_2), 2)$.
holdsAt$(coords(id_1, 201, 454), 2)$.
holdsAt$(coords(id_2, 227, 440), 2)$
holdsAt$(direction(id_1, 275), 2)$
holdsAt$(direction(id_2, 278), 2)$

Annotation for time 2:

holdsAt$(move(id_1, id_2), 2)$

(d)

Two Domain-specific axioms:

initiatedAt$(move(X, Y), T) \leftarrow$
  happensAt$(walk(X), T)$,
  happensAt$(walk(Y), T)$,
  $distLessThan(X, Y, 25, T)$,
  $dirLessThan(X, Y, 45, T)$

terminatedAt$(move(X, Y), T) \leftarrow$
  happensAt$(inactive(X), T)$,
  $distMoreThan(X, Y, 30, T)$

Table 1: **(a), (b)** The basic predicates and the domain-independent axioms of EC. **(c)** Example data from activity recognition. For example, at time point 1 person with id_1 is *walking*, her (X, Y) coordinates are $(201, 454)$ and her direction is 270°. The annotation for the same time point states that persons with id_1 and id_2 are not moving together, in contrast to the annotation for time point 2. **(d)** An example of two domain-specific axioms in the EC. E.g. the first clause dictates that *moving together* between two persons X and Y is initiated at time T if both X and Y are walking at time T, their euclidean distance is less than 25 pixel positions and their difference in direction is less than 45°. The second clause dictates that *moving together* between X and Y is terminated at time T if one of them is standing still at time T (exhibits an inactive behavior) and their euclidean distance at T is greater that 30.

As a running example we use the task of activity recognition, as defined in the CAVIAR project[4]. The CAVIAR dataset consists of 28 videos of actors performing a set of activities. Manual annotation (performed by the CAVIAR team) provides ground truth for two activity types. The first type corresponds to simple events and consists of the activities of a person at a certain video frame/time point, such as *walking*, or *standing still*. The second activity type corresponds to complex events and consists of activities that involve more than one person, e.g. two people *meeting each other*, or *moving together*. The goal is to recognize complex events as combinations of simple events and additional contextual knowledge, such as a person's direction and position.

Table 1(c) presents some example CAVIAR data, consisting of a narrative of simple events in terms of happensAt/2, expressing people's short-term activities, and context properties in terms of holdsAt/2, denoting people' coordinates and direction. Table 1(c) also shows the annotation of complex events (long-term activities) for each time-point in the narrative. Negated complex events' annotation is obtained via the closed-world

[4] http://homepages.inf.ed.ac.uk/rbf/CAVIARDATA1/

assumption (although both positive and negated annotation atoms are presented in Table 1(c), to avoid confusion). Table 1(d) presents two domain-specific axioms in the EC.

The learning task we address in this work is to learn definitions of complex events, in the form of domain-specific axioms in the EC, i.e. initiation and termination conditions of complex events, as in Table 1(d). The training data consists of a set of *Herbrand interpretations*, i.e. sets of narrative atoms, annotated by complex event instances, as in Table 1(c). Given such a training set $\mathcal{I}$ and some background knowledge B, the goal is to find a theory H that accounts for as many positive and as few negative examples as possible, throughout $\mathcal{I}$. Given an interpretation $I \in \mathcal{I}$, a positive (resp. negative) example is a complex event atom $\alpha \in I$ (resp. $\alpha \notin I$). We assume an online learning setting, in the sense that a learning algorithm is allowed only a single-pass over $\mathcal{I}$.

3.1 Online Learning of Markov Logic Networks with OSLα

OSLα [27] builds on the OSL [18] algorithm for online learning of Markov Logic Networks (MLN). An MLN is a set of weighted first-order logic rules. Along with a set of domain constants, it defines a ground Markov network containing one feature for each grounding of a rule in the MLN, with the corresponding weight. Learning an MLN consists of learning its structure (the rules in the MLN) and their weights.

OSLα works by constantly updating the rules in an MLN in the face of new interpretations that stream-in, by adding new rules and updating the weights of existing ones. At time t, OSL receives an interpretation I_t and uses its current hypothesis, H_t, to "make a prediction", i.e. to infer the truth values of *query atoms*, given the *evidence atoms* in I_t. In the learning problem that we address in this work, query atoms are instances of initiatedAt/2 and terminatedAt/2 predicates, the "target" predicates defining complex events, while evidence atoms are instances of predicates that declare simple events' occurrences, or other contextual knowledge, like happensAt/2 and $distLessThan/4$ respectively (see also Table 1(c)).

To infer the query atoms' truth values given the interpretation I_t, OSLα uses Maximum Aposteriori (MAP) inference [18], which amounts to finding the truth assignment to the query atoms that maximizes the sum of the weights of H_t's rules satisfied by I_t. This is a weighted MAX-SAT problem, whose solution OSLα efficiently approximates using LP-relaxed Integer Linear Programming, as in [17]. The inferred truth values of the query atoms, resulting from MAP inference may differ from the true ones, dictated by the annotation in I_t. The mistakes may be either False Negatives (*FNs*) or False Positives (*FPs*). Since *FNs* are query atoms which are not entailed by an existing rule in the inferred interpretation, in the next step OSLα constructs a set of new rules as a remedy. To this end, OSLα represents the current interpretation I_t as a hypergraph with the constants of I_t as the nodes and each true ground atom δ in I_t as a hyperedge connecting the nodes (constants) that appear as δ's arguments. Each inferred *FN* query atom α is used as a "seed" to generate new rules, by finding all paths in the hypergraph, up to a user-defined maximum length, that connect α's constants. Each such path consists of hyperedges, corresponding to a conjunction of true ground atoms connected by their arguments. The path is turned into a rule with this conjunction in the body and the seed *FN* atom in the head, and a lifted version of this rule is obtained by variabilizing the

ground arguments. By construction, each such rule logically entails the seed *FN* atom it was generated from.

This technique of *relational pathfinding* may generate a large number of rules from each inferred *FN* query atom, and many of these rules are not useful. Thus, OSLα relies on L_1-regularized weight learning, which in the long run, pushes the weights of non-useful rules to zero. Weight learning is also OSLα's way to handle *FP* query atoms in the inferred state, which are due to erroneously satisfied rules in the current theory H_t. These rules are penalized by reducing their weights. OSLα uses an AdaGrad-based [13] weight learning technique, which supports L_1-regularization.

AdaGrad is a subgradient-based method for online convex optimization, i.e. at each step it updates a feature vector, based on the subgradient of a convex loss function of the features. Seeing the rules in an MLN theory $H = \{r_1, \ldots, r_n\}$ as the feature vector that needs to be optimized, the authors in [18] use a simple variant of the hinge-loss as a loss function, whose subgradient is the vector $-\langle \Delta g_1, \ldots \Delta g_n \rangle$, where Δg_i denotes the difference between the true groundings of the i-th rule in the actual true state and the MAP-inferred state respectively. Based on this difference, AdaGrad updates the weight w_i^t of the i-th rule in the theory at time t by:

$$w_i^{t+1} = sign(w_i^t - \frac{\eta}{C_i^t} \Delta g_i^t)\ max\{0, |w_i^t - \frac{\eta}{C_t^i} \Delta g_i^t| - \lambda \frac{\eta}{C_t^i}\} \tag{1}$$

where t-superscripts in terms denote the respective values at time t, η is a learning rate parameter, λ is a regularization parameter and $C_i^t = \delta + \sqrt{\sum_{j=1}^{t} (\Delta g_i^j)^2}$ is a term that expresses the rule's quality so far, as reflected by the accumulated sum of Δg_i's, amounting to the i-th rule's past mistakes (plus a $\delta \geq 0$ to avoid division by zero in η / C_i^t). The C_i^t term gives an adaptive flavour to the algorithm, since the magnitude of a weight update via the term $|w_i^t - \frac{\eta}{C_t^i} \Delta g_i^t|$ in Equation (1), is affected by the rule's previous history, in addition to its current mistakes, expressed by Δg_i^t. The regularization term in Equation (1), $\lambda \frac{\eta}{C_t^i}$, is the amount by which the i-th rule's weight is discounted when $\Delta g_i^t = 0$. This is to eventually push to zero the weights of irrelevant rules, which have very few, or even no groundings in the training interpretations.

In contrast to its predecessor, OSL, OSLα uses specialized techniques for pruning large parts of the hypergraph structure to enhance efficiency when learning with the EC. However, its rule generation technique remains an important bottleneck. Repeatedly searching for paths in the hypergraph structure is expensive in its own right, but it also tends to blindly generate large sets of rules, which, in turn, increases the cost of MAP inference during learning.

4 Learning Weighted Rules with WoLED

Aiming to improve the efficiency of online learning with the EC in MLN, we propose WoLED (Weighted Online Learning of Event Definitions), an extension of the OLED crisp online ILP learner [21], to an MLN setting. OLED draws inspiration from the VFDT (Very Fast Decision Trees) algorithm [10], whose online strategy is based on the Hoeffding bound [16]. Given a random variable X with range in $[0, 1]$ and an observed mean $\overline{X}$ of its values after n independent observations, the Hoeffding Bound states that,

with probability $1-\delta$, the true mean $\hat{X}$ of the variable lies in an interval $(\overline{X}-\epsilon, \overline{X}+\epsilon)$, where $\epsilon = \sqrt{\frac{ln(1/\delta)}{2n}}$.

OLED learns a rule r with a hill-climbing process, where literals are gradually added to r's body yielding *specializations* of r of progressively higher quality, which is assessed by some scoring function G, based on the positive and negative examples that a rule entails. This strategy is common in ILP, where at each specialization step (addition of a literal), a number of candidate specializations of the parent rule are evaluated on the entire training set and the best one is selected. OLED adapts this strategy to an online setting, using the Hoeffding bound, as follows: Assume that after having evaluated r and a number of its candidate specializations on n examples, r_1 is r's specialization with the highest mean G-score $\overline{G}$ and r_2 is the second-best one, i.e. $\Delta\overline{G} = \overline{G}(r_1) - \overline{G}(r_2) > 0$. Then by the Hoeffding bound we have that for the true mean of the scores' difference $\Delta\hat{G}$ it holds that $\Delta\hat{G} > \Delta\overline{G} - \epsilon$, with probability $1-\delta$, where $\epsilon = \sqrt{\frac{ln(1/\delta)}{2n}}$. Hence, if $\Delta\overline{G} > \epsilon$, then $\Delta\hat{G} > 0$, implying that r_1 is indeed the best specialization, with probability $1-\delta$. In order to decide which specialization to select, it thus suffices to evaluate the specializations on examples from the stream until $\Delta\overline{G} > \epsilon$. Each of these examples is processed once, thus giving rise to a single-pass rule learning strategy.

Positive (resp. negative) examples in this setting are true (resp. false) instances of target predicates, which are present in (resp. generated via the closed-world assumption from) the incoming training interpretations. Such instances correspond to query atoms in an MLN setting (see Section 3.1), therefore we henceforth refer to positive and negative examples as true and false query atoms respectively.

The literals used for specialization are drawn from a *bottom rule* [8], denoted by $\perp$, a most-constrained rule that entails a single true query atom. A bottom rule is usually too restrictive to be used for classification and its purpose is to define a space of potentially good rules, consisting of those that θ-subsume $\perp$. OLED moves through this search space in a top-down fashion, starting from an empty-bodied rule $head(\perp) \leftarrow true$ and considering at each time specializations that result by the addition of a fixed number of literals from $\perp$ to the body of a parent rule (the default is one literal).

4.1 WoLED

Our approach is based on combining OLED's rule learning technique with OSLα's weight learning strategy. Contrary to OSLα's approach, where once a rule is generated, its structure remains fixed and the only way to improve its quality is by tuning its weight, WoLED progressively learns both the structure of a rule and its weight, by jointly optimizing them together. To this end, a rule is gradually specialized via the online hill-climbing strategy described earlier, while constantly updating the weight of each such specialization. Given a rule r and a set of r's specializations S_r, WoLED learns weights for r and each $r' \in S_r$, and uses Hoeffding tests to identify over time, a rule $r_1 \in S_r$ of higher quality, as assessed by a combination of r_1's structure and weight. Once that happens, r is replaced by r_1 and the process continues for as long as new specializations of r_1 improve its quality. To learn weights, WoLED replaces OLED's crisp logical inference with MAP inference and uses OSLα's mistake-driven

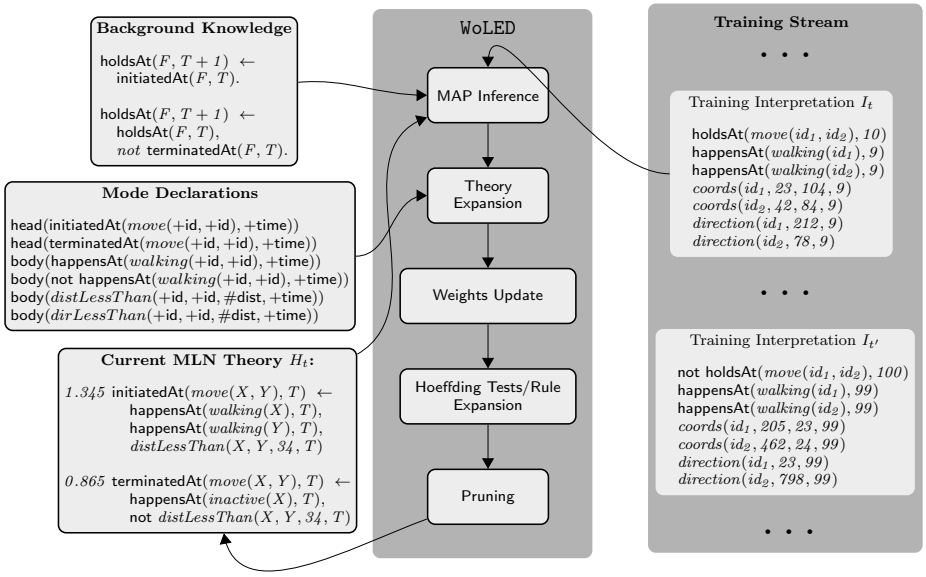


Fig. 1: Illustration of WoLED's high-level strategy.

weight learning technique described in Section 3.1, which updates a rule's weight based on the query atoms that the rule misclassifies in the MAP-inferred state.

WoLED's high-level strategy is illustrated in Figure 1. At each point in time WoLED maintains a theory $H_t = \{r_1, \ldots, r_n\}$, and for each rule r_i, a set of current specializations $\{r_i^1, \ldots, r_i^n\}$ and an associated bottom rule $\bot_{r_i}$. At time t, WoLED receives the t-th training interpretation I_t and it subsequently goes through a five-step process. In the first step it performs MAP inference on I_t, using the current theory H_t and the background knowledge. The MAP-inferred interpretation is checked for *FN* query atoms. If such an atom α exists, meaning that no existing rule in H_t entails it, WoLED proceeds to a "theory expansion" step, where it starts growing a new rule r_{i+1} to account for that. This amounts to using the *FN* atom α as a "seed" to generate a bottom rule $\bot_{r_{i+1}}$ and adding to H the empty-bodied rule $head(\bot_{r_{i+1}}) \leftarrow true$. From that point on this rule is gradually specialized using literals from $\bot_{r_{i+1}}$. The generation of $\bot_{r_{i+1}}$ is guided by a set of *mode declarations* [8] (see Figure 1), a form of language bias specifying the signatures of literals that are allowed to be placed in the head or the body of a rule, in addition to the types of their arguments (e.g. time, or id in the declarations of Figure 1) and whether they correspond to variables or constants (indicated respectively by "+" and "#" in the declarations of Figure 1).

After a potential theory expansion a weight update step follows, where the weights of each rule $r \in H_t$ and the weights of each one of its current specializations are updated, based on their mistakes on the MAP-inferred state generated previously. A Hoeffding test for each rule $r \in H_t$ follows, and if a test succeeds for some rule r, it is "expanded", i.e. r's structure and weight are replaced by the ones of its best-scoring specialization, r_1. The final step is responsible for pruning the current theory, i.e. re-

Algorithm 1 WoLED

Input: $\mathcal{I}$: A stream of training interpretations; M: Mode declarations; B: Background knowledge; G: Rule evaluation function; δ_h: Confidence for the Hoeffding test; η, λ, δ_a: AdaGrad's parameters; d: Specialization depth; N_{min}: Warm-up period; $Score_{min}$: G-score quality threshold; H: A (possibly empty) set of first order rules.

1: **for each** $I \in \mathcal{I}$ **do**
2: $H_{active} = \{r \in H \ : r$ has been evaluated on at least N_{min} examples$\}$
3: $I_{MAP} :=$ the MAP-inferred interpetation of $H_{active} \cup B$ on I.
4: **for each** true query atom $\alpha \in I \setminus I_{MAP}$ **do**
5: Generate a bottom rule $\perp$ from $I \cup B$, using the mode declarations and α as a seed atom.
6: $r := head(\perp) \leftarrow true$
7: $TPs(r), FPs(r), FNs(r), Subgradients(r) := 0$; $\perp_r := \perp$
8: $H \leftarrow H \cup \{r\}$
9: **for each** $r \in H$ **do**
10: $\rho_d(r) := \{head(r) \leftarrow body(r) \wedge D \mid D \subset body(\perp_r)$ and $|D| \leq d\}$
11: `UpdateWeights`(r, I_{MAP})
12: **for each** $r' \in \rho_d(r)$ **do**
13: `UpdateWeights`(r', I_{MAP})
14: $\Delta\bar{G} := \bar{G}(r_1) - \bar{G}(r_2)$, where $r_1, r_2 \in \rho_d(r)$ are r's best and second-best specializations.
15: $\epsilon := \sqrt{\frac{ln(1/\delta_h)}{2N_r}}$, where N_r is the sum of r's groundings so far.
16: $\bar{\epsilon} :=$ the mean value of $\epsilon = \sqrt{\frac{ln(1/\delta_h)}{2N_{r_k}}}$ observed so far any rule r_k.
17: **if** $[\Delta\bar{G} > \epsilon$ **or** $\Delta\bar{G} < \epsilon < \bar{\epsilon}]$ **and** $\bar{G}(r_1) > \bar{G}(r)$ **then**
18: Replace r by r_1.
19: $\bar{N_s} :=$ the average number of $\mathcal{O}(\frac{1}{\epsilon^2} ln \frac{1}{\delta_h})$ examples for which the Hoeffding test has succeeded so far during the learning process.
20: **if** r is unchanged for a period longer than $\bar{N_s}$ and $Score_{min} - \bar{G}(r) > \epsilon$ **then**
21: Remove r from H.
22: **return** H

moving low-quality rules. The current version of the weighted theory is output and the training loop continues (see Figure 1).

We next go into some details of WoLED's functionality, using the pseudocode in Algorithm 1. The input to Algorithm 1 consists of the background knowledge, a rule-scoring function G, based on the true positive (*TP*), false positive (*FP*) and false negative (*FN*) examples entailed by a rule, AdaGrad's parameters η, δ_a, λ, discussed in Section 3.1, the confidence parameter for the Hoeffding test δ_h and a set of mode declarations. Additionally, a minimum G-score value for acceptable rules, a "warm-up" parameter N_{min} and a "specialization depth" parameter d, to be explained shortly.

The MAP-inferred state is generated in lines 2–3 of Algorithm 1, using the *active fragment* of the theory. The latter consists of those rules in the theory that have been evaluated on at least N_{min} examples (line 2), where N_{min} is the input "warm-up" parameter. This is to avoid using rules which are too premature and useless for inference, such as an empty-bodied rule that has just been created.

Algorithm 2 `UpdateWeights`$(r, \eta, \delta_a, \lambda, I_{MAP}, I_{TRUE})$:
Input: r: a rule; η, λ, δ_a: AdaGrad's learning rate, regularization parameter and smoothness parameter respectively; I_{MAP}, I_{TRUE}: the MAP-inferred state and the true state respectively for a training interpretation I.

1: $w_r :=$ the weight of rule r.
2: $\Delta g_r :=$ the difference in true groundings of rule r in the inferred state I_{MAP} and the true state I_{TRUE}.
3: $Subgradients(r) \leftarrow Subgradients(r) + (\Delta g_r)^2$.
4: $C_r := \delta_\alpha + \sqrt{Subgradients(r)}$
5: $w_r \leftarrow sign(w_r - \frac{\eta}{C_r}\Delta g_r)\ max\{0, |w_r - \frac{\eta}{C_r}\Delta g_r| - \lambda\frac{\eta}{C_r}\}$
6: $TPs(r) \leftarrow TPs(r) + |\{\alpha \in I_{MAP} \cap I_{TRUE} : \alpha \text{ is a grounding of } head(r)\}|$.
7: $FPs(r) \leftarrow FPs(r) + |\{\alpha \in I_{MAP} \setminus I_{TRUE} : \alpha \text{ is a grounding of } head(r)\}|$.
8: $FNs(r) \leftarrow FNs(r) + |\{\alpha \in I_{TRUE} \setminus I_{MAP} : \alpha \text{ is a grounding of } head(r)\}|$.

In lines 4-8 of Algorithm 1, new rules are generated for each *FN* atom in the MAP-inferred state, as described earlier. In addition to its corresponding bottom rule to draw literals for specialization, each rule is also equipped with a number of counters, which accumulate the *TP*, *FP* and *FN* instances entailed by the rule over time, to use for calculating its G-score; and an accumulator, denoted by $Subgradients(r)$, which is meant to store the history (sum) of the rule's mistakes throughout the learning process (see the term C_i^t in Equation (1)). As explained in Section 3.1, a rule's mistakes w.r.t. each incoming interpretation is a coordinate in the the subgradient vector of AdaGrad's loss function (hence the name "$Subgradients(r)$" for their accumulator) and they affect the magnitude of a rule's weight update. Therefore, $Subgradients(r)$ is used for updating r's weight with AdaGrad.

Updating the weights of each rule $r \in H$, as well as the weights of r's candidate specializations follows, in lines 9-13 of Algorithm 1 (r's specializations are denoted by $\rho_d(r)$ (line 10), where d is the specialization depth parameter mentioned earlier, controlling the "depth" of allowed specializations, which are considered at each time). The weight-update process is presented in Algorithm 2. It uses AdaGrad's strategy discussed in Section 3.1. The difference between r's true groundings in the true state and the MAP-inferred one is first calculated (line 2), it's square is added to r's $Subgradients$ accumulator (line 3 – see the C_i^t term in Equation (1), Section 3.1) and then r's weight is updated (line 5). Algorithm 2 is also responsible for updating the TP, FP, FN counters for a rule r, which are used to calculate its G-score.

The Hoeffding test follows to decide if a rule should be specialized (lines 14-18, Algorithm 1). A rule is specialized either if the test succeeds ($\Delta\bar{G} > \epsilon$), or if a tie-breaking condition is met ($\Delta\bar{G} < \epsilon < \tau$), where τ is a threshold set to the mean value of ϵ observed so far. Also, to ensure that no rule r is replaced by a specialization of lower quality, we demand that $\bar{G}(r_1) > \bar{G}(r)$, where r_1 is r's best-scoring specialization indicated by the Hoeffding test.

The final step is responsible for removing rules of low quality (lines 19-21, Algorithm 1). A rule is removed if it remains unchanged (is not specialized) for a significantly large period of time, set to the average number of $\mathcal{O}(\frac{1}{\epsilon^2} ln \frac{1}{\delta_h})$ examples for which the Hoeffding test has succeeded so far, and there is enough confidence, via an

additional Hoeffding test, that its mean G-score is lower than a minimum acceptable G-score $Score_{min}$.

5 Experimental Evaluation

We present an experimental evaluation of WoLED on CAVIAR, a benchmark dataset for activity recognition (see Section 3 for CAVIAR's description). All experiments were conducted on Debian Linux machine with a 3.6GHz processor and 16GB of RAM. The code and the data to reproduce the experiments are available online[5]. WoLED is implemented in Scala, using the Clingo answer set solver[6] for grounding and the lpsolve[7] Linear Programming solver for probabilistic MAP-inference, on top of the LoMRF[8] platform, an implementation of MLN. The OSLα version to which we compare in these experiments also relies on lpsolve, but uses LoMRF's custom grounder.

5.1 Comparison with Related Online and Batch Learners

In our first experiment we compare WoLED with (i) OSLα and OSL [18], discussed in Section 3.1; (ii) The crisp version of OLED; (iii) EC_{crisp}, a hand-crafted set of crisp rules for CAVIAR; (iv) MaxMargin, an MLN consisting of EC_{crisp}'s rules, with weights optimized by the the Max-Margin weight learning method of [17]; (v) XHAIL, a batch, crisp ILP learner using a combination of inductive and adbuctive logic programming.The rules used by EC_{crisp} and MaxMargin may be found in [30].

MaxMargin was selected because it was shown to achieve good results on CAVIAR [30], while XHAIL was selected because it is one of the few existing ILP algorithms capable of learning theories in the EC.

A fragment of the CAVIAR dataset has been used in previous work to evaluate OSLα and MaxMargin's performance [27, 30]. To compare to these approaches we therefore used this fragment in our first experiment. The target complex events in this dataset are related to two persons *meeting each other* or *moving together* and the training data consist of the parts of CAVIAR where these complex events occur. The fragment dataset contains a total of 25,738 training interpretations. The results with OLED were achieved using a significance parameter $\delta_h = 10^{-2}$ for the Hoeffding test, a rule pruning threshold $Score_{min}$ (see also Algorithm 1) of 0.8 for *meeting* and 0.7 for *moving* and a warm-up parameter of $N_{min} = 1000$ examples. WoLED also used this parameter configuration, in addition to $\eta = 1.0, \lambda = 0.01, \delta_\alpha = 1.0$ for weight learning with AdaGrad. These parameters were reported in [27] and were also used with OSLα/OSL.

The results were obtained using 10-fold cross validation and are presented in Table 2(a) in the form of *precision, recall* and f_1*-score*. These statistics were micro-averaged over the instances of recognized complex events from each fold.Table 2(a) also presents average training times per fold for all approaches except EC_{crisp}, where there is no training involved, average theory sizes for OLED, OSLα, and XHAIL, as well as the

[5] https://github.com/nkatzz/OLED
[6] https://potassco.org/clingo/
[7] https://sourceforge.net/projects/lpsolve/
[8] https://github.com/anskarl/LoMRF

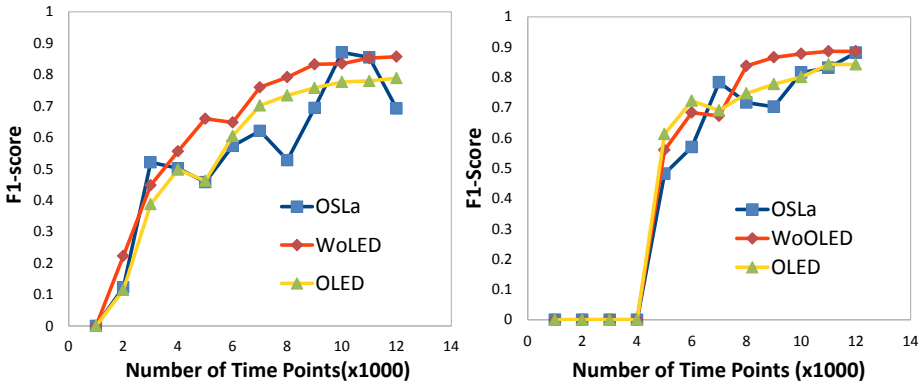
		Method	Precision	Recall	F_1-score	Theory size	Time (sec)
(a)	*Moving*	$\mathsf{EC}_{\mathsf{crisp}}$	**0.909**	0.634	0.751	28	–
		OLED	0.867	0.724	0.789	34	**28**
		WoLED	0.882	0.835	0.857	30	59
		OSLα	0.837	0.590	0.692	3316	1300
		OSL	-	-	-	-	$>$ 25 hrs
		MaxMargin	0.844	**0.941**	**0.890**	28	1692
		XHAIL	0.779	0.914	0.841	**14**	7836
	Meeting	$\mathsf{EC}_{\mathsf{crisp}}$	0.687	0.855	0.762	23	–
		OLED	**0.947**	0.760	0.843	31	**22**
		WoLED	0.892	0.888	**0.889**	29	52
		OSLα	0.902	0.863	0.882	1231	180
		OSL	-	-	-	-	$>$ 25 hrs
		MaxMargin	0.919	0.813	0.863	23	1133
		XHAIL	0.804	**0.927**	0.861	**15**	7248
(b)	*Moving*	OLED	0.682	0.787	0.730	**38**	**63**
		WoLED	**0.783**	**0.821**	**0.801**	51	108
		$\mathsf{EC}_{\mathsf{crisp}}$	0.721	0.639	0.677	28	–
	Meeting	OLED	0.701	0.886	0.782	**41**	**43**
		WoLED	**0.808**	**0.877**	**0.841**	56	98
		$\mathsf{EC}_{\mathsf{crisp}}$	0.644	0.855	0.735	23	–

Table 2: Experimental results on (a) the CAVIAR fragment of [30] (top) and (b) the complete CAVIAR dataset (bottom).

fixed theory size of $\mathsf{EC}_{\mathsf{crisp}}$ and MaxMargin. The reported theory sizes are in the form of total number of literals in a theory. The online methods were allowed only a single-pass over the training data.

WoLED achieves the best F_1-score for *meeting* and the second-best F_1-score for *moving*, right after the batch weight optimizer MaxMargin. This is a notable result. Moreover, this gain in predictive accuracy comes with a tolerable decrease in efficiency of approximately half a minute, as compared to OLED's training times, which are the best among all learners. This extra overhead in training time for WoLED is due to the cost of the probabilistic MAP-inference, which replaces OLED's crisp logical inference to allow for weight optimization. Regarding theory sizes, WoLED outputs hypotheses comparable in size with the hand-crafted knowledge base, and much more compressed as opposed to OSLα. This is another notable result. XHAIL learns the most compressed hypotheses, since it is a batch learner, which also explains its increased training times.MaxMargin, also has high training times, paying the price of batch (weight) optimization.

OSL was unable to process the dataset within 25 hours, at which time training was terminated. The reasons for that are related to it being unable to take advantage of the background knowledge, thus it is practically unable to learn with the Event Calculus



(a) *moving* (b) *meeting*

Fig. 2: Online holdout evaluation on CAVIAR.

[27]. OSLα overcomes OSL's difficulties, but learns unnecessarily large theories, which differ in size by several orders of magnitude from all others learners'. In turn, this affects OSLα's training times, which are also increased. In contrast, WoLED achieves improved predictive accuracy and compressed theories with minimal training overhead.

Figure 2 presents the holdout evaluation [15] for the online learners compared in this experiment. Holdout evaluation consists of assessing the quality of an online learner on a holdout test set, at regular time intervals during learning, thus obtaining a learning curve of its performance over time. Figure 2 presents average F_1-scores, obtained by performing holdout evaluation on each fold of the tenfold cross-validation process: at each fold, each learner's theory is evaluated on the fold's test set every 1000 time points and the F_1-scores from each evaluation point are averaged over all ten folds.

WoLED and OLED have an adequate performance, with relatively smooth learning curves, while they eventually converge to stable theories of acceptable performance. Moreover, WoLED outperforms both OLED and OSLα in most of the evaluation process.

In contrast to the online behaviour of WoLED and OLED, OSLα's performance exhibits abrupt fluctuations. For *moving* in particular, OSLα's average F_1-score reaches its peak (0.87) after processing data from approximately 10,000 time points, and then it drops significantly until the final average F_1-score value of 0.69 reported in Table 2. This behavior may be attributed to OSLα's rule generation strategy. Contrary to WoLED, which uses Hoeffding tests to select rules with significant heuristic value, OSLα greedily adds new rules to the current theory, so as to locally improve its performance, without taking into account the new rules' quality on larger portions of the data. Overall, this results in poor online performance, since rules with no quality guarantees on the training set may be responsible for a large number of mistakes on unseen data, by e.g. fitting the noise in the training data. OSLα relies solely on weight learning to minimize the weights of low-quality rules in the long run. However, OSLα's holdout evaluation indicates that in principle this requires larger training sets, since, at least in the case of *moving*, OSLα's theories exhibit no sign of convergence. On the other hand, OSLα's in-

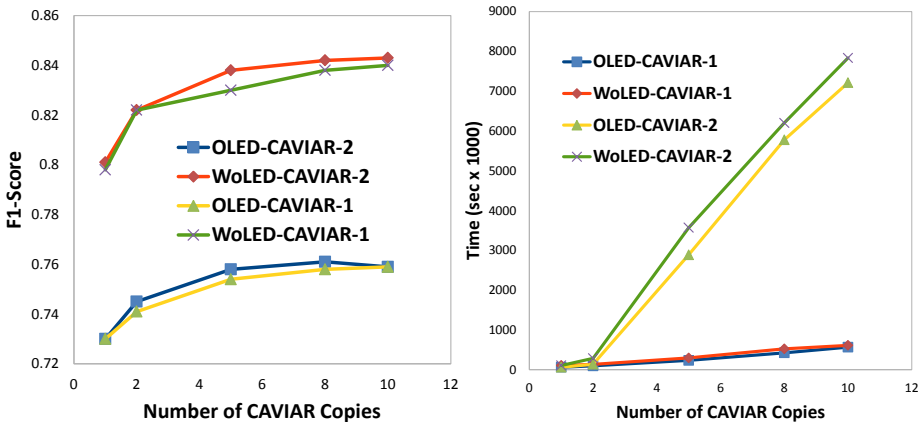


Fig. 3: Evaluation on larger data volumes for the *moving* complex event.

creased training times reported in Table 2, due to the ever-increasing cost of maintaining unnecessarily large theories, indicate that training on larger datasets is impractical.

5.2 Evaluation on Larger Data Volumes

In this section we evaluate WoLED on larger data volumes, starting with the entire CAVIAR dataset, which consists of 282,067 interpretations, in contrast to 25,738 interpretations in the CAVIAR fragment. Due to the increased training times of OSLα, XHAIL and MaxMargin, we did not experiment with these algorithms. The target complex events were *meeting* and *moving* as previously. The additional training data (i.e. those not contained in the CAVIAR fragment) were negative instances for both complex events (recall that the parts of CAVIAR where *meeting* and *moving* occur were already contained in the CAVIAR fragment). This way, the dataset used in this experiment is much more imbalanced than the one used in the previous experiment. The parameter configuration for the two learners was as reported in Section 5.1. The results were obtained via tenfold cross-validation and are presented in Table 2(b).

The average F_1-score for both algorithms is decreased, as compared to the previous experiment, due to the increased number of false positives, caused by the large number of additional negative instances. WoLED significantly outperforms OLED for both complex events, at the price of a tolerable increase in training times.

To test our approach further we used larger datasets generated from CAVIAR in two different settings. In the first setting, to which we henceforth refer as CAVIAR-1, we generated datasets by sequentially appending copies of the original CAVIAR dataset, incrementing the time-stamps in each copy. Therefore, training with datasets in the CAVIAR-1 setting amounts to re-iterating over the original dataset a number of times. In the second setting, to which we refer as CAVIAR-2, datasets were also obtained from copies of CAVIAR, but this time the time-stamps in the data were left intact and each copy differed from the others in the constants referring to the tracked entities (persons,

objects) that appear in simple and complex events. In each copy of the dataset, the coordinates of each entity p differ by a fixed offset from the coordinates of the entity of the original dataset that p mirrors. The copies were "merged", grouping together by time-stamp the data from each copy. Therefore, the number of constants in each training interpretation in datasets of the CAVIAR-2 setting is multiplied by the number of copies used to generate the dataset. We performed experiments on datasets obtained from 2, 5, 8 and 10 CAVIAR copies for both settings. The target concept was *moving* and on each dataset we used tenfold cross-validation to measure F_1-scores and training times.

The results are presented in Figure 3. F_1-scores improve with larger data volumes for both learners, slightly more so with the datasets in the CAVIAR-2 setting, while WoLED achieves better F_1-scores than OLED thanks to weight learning. Training times increase slowly with the data size in the "easier" CAVIAR-1 setting, where both learners require 8-10 minutes on average to learn from the largest dataset in this setting. In contrast, training times increase abruptly in the harder CAVIAR-2 setting, where learning from the largest dataset requires more than 2.5 hours on average for both learners. This is due to the additional domain constants in the datasets of the CAVIAR-2 setting, which result in exponentially larger ground theories.

6 Conclusions and Future Work

We presented an algorithm for online learning of event definitions in the form of Event Calculus theories in the MLN semantics. We extended an online ILP algorithm to a statistical relational learning setting via online weight optimization. We evaluated our approach on an activity recognition application, showing that it outperforms both its crisp predecessor and competing algorithms for online learning in MLN. There are several directions for further work. We plan to improve scalability using parallel/distributed learning, along the lines of [22]. We also plan to evaluate different algorithms for online weight optimization and develop methodologies for online hyper-parameter adaptation.

References

1. E. Alevizos, A. Skarlatidis, A. Artikis, and G. Paliouras. Probabilistic complex event recognition: A survey. *ACM Computing Surveys*, to appear, 2018.
2. A. Artikis, M. Sergot, and G. Paliouras. An event calculus for event recognition. *Knowledge and Data Engineering, IEEE Transactions on*, 27(4):895–908, 2015.
3. A. Artikis, A. Skarlatidis, F. Portet, and G. Paliouras. Logic-based event recognition. *The Knowledge Engineering Review*, 27(4):469–506, 2012.
4. D. Athakravi, D. Corapi, K. Broda, and A. Russo. Learning through hypothesis refinement using answer set programming. In *ILP-2013*, pages 31–46. Springer, 2013.
5. D. Corapi, A. Russo, and E. Lupu. Inductive logic programming as abductive search. In *ICLP-2010*, pages 54–63, 2010.
6. D. Corapi, D. Sykes, K. Inoue, and A. Russo. Probabilistic rule learning in nonmonotonic domains. In *International Workshop on Computational Logic in Multi-Agent Systems*, pages 243–258. Springer, 2011.
7. G. Cugola and A. Margara. Processing flows of information: From data stream to complex event processing. *ACM Computing Surveys (CSUR)*, 44(3):15, 2012.

8. L. De Raedt. *Logical and relational learning*. Springer Science & Business Media, 2008.
9. L. De Raedt, K. Kersting, S. Natarajan, and D. Poole. Statistical relational artificial intelligence: Logic, probability, and computation. *Synthesis Lectures on Artificial Intelligence and Machine Learning*, 10(2):1–189, 2016.
10. P. Domingos and G. Hulten. Mining high-speed data streams. In *ACM SIGKDD*, pages 71–80. ACM, 2000.
11. S. Dragiev, A. Russo, K. Broda, M. Law, and C. Turliuc. An abductive-inductive algorithm for probabilistic inductive logic programming. In *Proceedings of the 26th International Conference on Inductive Logic Programming (Short papers), London, UK, 2016.*, pages 20–26, 2016.
12. A. Dries and L. De Raedt. Towards clausal discovery for stream mining. In *ILP-2009*, pages 9–16. Springer, 2009.
13. J. Duchi, E. Hazan, and Y. Singer. Adaptive subgradient methods for online learning and stochastic optimization. *Journal of Machine Learning Research*, 12(Jul):2121–2159, 2011.
14. J. Gama. *Knowledge discovery from data streams*. CRC Press, 2010.
15. J. Gama, R. Sebastião, and P. P. Rodrigues. On evaluating stream learning algorithms. *Machine learning*, 90(3):317–346, 2013.
16. W. Hoeffding. Probability inequalities for sums of bounded random variables. *Journal of the American statistical association*, 58(301):13–30, 1963.
17. T. N. Huynh and R. J. Mooney. Max-margin weight learning for markov logic networks. In *Joint European Conference on Machine Learning and Knowledge Discovery in Databases*, pages 564–579. Springer, 2009.
18. T. N. Huynh and R. J. Mooney. Online structure learning for markov logic networks. In *ECML-2011*, pages 81–96. Springer, 2011.
19. N. Katzouris. Scalable relational learning for event recognition. *PhD Thesis, University of Athens*, http://users.iit.demokritos.gr/ nkatz/papers/nkatz-phd.pdf, 2017.
20. N. Katzouris, A. Artikis, and G. Paliouras. Incremental learning of event definitions with inductive logic programming. *Machine Learning*, 100(2-3):555–585, 2015.
21. N. Katzouris, A. Artikis, and G. Paliouras. Online learning of event definitions. *Theory and Practice of Logic Programming*, 16(5-6):817–833, 2016.
22. N. Katzouris, A. Artikis, and G. Paliouras. Parallel online learning of event definitions. In *ILP*, 2017.
23. R. Kowalski and M. Sergot. A logic-based calculus of events. *New Generation Computing*, 4(1):67–95, 1986.
24. M. Law, A. Russo, and K. Broda. Iterative learning of answer set programs from context dependent examples. *Theory and Practice of Logic Programming*, 16(5-6):834–848, 2016.
25. N. Littlestone. Learning quickly when irrelevant attributes abound: A new linear-threshold algorithm. *Machine learning*, 2(4):285–318, 1988.
26. A. Margara, G. Cugola, and G. Tamburrelli. Learning from the past: automated rule generation for complex event processing. In *Proceedings of the 8th ACM International Conference on Distributed Event-Based Systems*, pages 47–58. ACM, 2014.
27. E. Michelioudakis, A. Skarlatidis, G. Paliouras, and A. Artikis. Osla: Online structure learning using background knowledge axiomatization. In *ECML*, pages 232–247. Springer, 2016.
28. O. Ray. Nonmonotonic abductive inductive learning. *J. Applied Logic*, 7(3):329–340, 2009.
29. M. Richardson and P. Domingos. Markov logic networks. *Machine learning*, 62(1-2):107–136, 2006.
30. A. Skarlatidis, G. Paliouras, A. Artikis, and G. Vouros. Probabilistic event calculus for event recognition. *ACM Transactions on Computational Logic (TOCL)*, 16(2):11, 2015.
31. A. Srinivasan and M. Bain. An empirical study of on-line models for relational data streams. *Machine Learning*, 106(2):243–276, 2017.