Learning to Search

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Learning to Search

structured prediction as search problem

A structured learning problem consists of

- an input space ${\mathcal X}$
- an output space \mathcal{Y}
- a fixed but unknown data distribution \mathcal{D} over $\mathcal{X} \times \mathcal{Y}$
- a loss function $\ell(y^*, \hat{y}) \to \mathbb{R}^+$, which measures the distance between the true (y^*) and predicted (\hat{y}) outputs.

The goal of structured learning is to use N samples, $\{x_i, y_i\}_{i=1}^N \sim \mathcal{D}$ and learn a mapping $f : \mathcal{X} \to \mathcal{Y}$, that minimizes the expected loss $\mathbb{E}_{(x,y)\sim \mathcal{D}}[\ell(y, f(x))]$



- the 0/1 loss, $\mathbb{I}[y \cdot w^{\top} x \leq 0]$, is unfortunately non-convex
- we already know that we should build some convex surrogate:

$$\ell(w, x) = \max(0, 1 - y \cdot w^{\top} x)$$

- this is called 'hinge-loss'
- it's convex and upper-bounds the 0/1 loss, $\ell_t(w, x) \ge \mathbb{I}[y \cdot w^\top x \le 0], \forall w$



- **1** receive x_t
- **2** predict $p_t = sign(w_t^\top x_t)$
- 3 suffer loss $\ell_t(w, x_t) = \max(0, 1 y \cdot w^\top x_t)$
- 4 update w_t

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We know already Follow-The-Regularized-Leader as a way to no-regret:

$$w_t = \arg\min\sum_{i=1}^t \ell_t(w, x_i) + R(w)$$

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We know already Follow-The-Regularized-Leader as a way to no-regret:

$$w_t = \arg\min\sum_{i=1}^t \ell_t(w, x_i) + \frac{1}{2\nu} ||w||_2^2$$

We also know that for convex functions it is sufficient to work with linearized losses, setting $z_t = \nabla_w \ell_t(w)$:

Theorem

Consider FTRL, linear losses $\ell_t(w) = w^{\top} z_t$, and regularization $R(w) = \frac{1}{2\nu} ||w||_2^2$ and $w, u \in S = \mathbb{R}^d$, then

$$R_T(u) \le \frac{1}{2\nu} ||u||_2^2 + \frac{\nu}{2} \sum_{t=1}^T ||z_t||_2^2.$$

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- when there is no error the gradient is zero
- \bullet z_t is the gradient

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- when there is no error the gradient is zero
- z_t is the gradient
- gradient of $w^{\top}x_t$ is just x_t

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• and let $R = \max ||x_t||$ (bounded features)

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- use the definition of regret and the surrogate upper-bounding property

$$\#errors \le \sum_{t} \ell(w_{t}) - \sum_{t} \ell(u) \le \frac{1}{2\nu} ||u||_{2}^{2} + \frac{\nu}{2} R^{2} \cdot \#errors$$

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$$\begin{split} \#errors &\leq \sum_{t} \ell(w_{t}) - \sum_{t} \ell(u) \leq \frac{1}{2\nu} ||u||_{2}^{2} + \frac{\nu}{2} R^{2} \cdot \#errors \\ \bullet \quad \text{set } \nu &= \frac{||u||}{R\sqrt{\#errors}} \text{ and rearrange} \\ \#errors - R ||u|| \sqrt{\#errors} - \sum_{t=1}^{T} \ell_{t}(u) \leq 0 \end{split}$$

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• let's assume realizability, $\exists u^*$, s.t. $y_t = sign((u^*)^\top x_t), \forall t$ $\# errors - R ||u^*|| \sqrt{\# errors} \leq 0$

if we additionaly assume, that $||u^*|| \leq 1/\gamma$ we get $\# errors \leq R^2/\gamma^2$

[Novikoff, 1962]

Structured Perceptron

Structured Perceptron [Collins and Roark, 2004]

- 1: **input:** training data $\{(x_t, y_t)\}_{t=1}^N$
- 2: **init**: $w_0 = 0$
- 3: for $t = 0, \ldots$ do
- 4: observe x_t, y_t
- 5: predict $\hat{y}_t = \arg \max_{y \in \mathcal{Y}(x)} w_t^\top \phi(x_t, y)$
- 6: **if** $\hat{y}_t \neq y_t$ then
- 7: update $w_{t+1} = w_t + \phi(x_t, y_t) \phi(x_t, \hat{y}_t)$

8: end if

Thm

$$\begin{split} & \text{If } ||\phi(x_t,y_t)|| \leq R \text{ and } \exists u, ||u|| = 1 \text{ such that } \forall t,y \in \mathcal{Y}(x_t): \\ & u^\top \phi(x_t,y_t) \geq u^\top \phi(x_t,y) + \gamma \text{, then} \end{split}$$

$$\# \text{ errors} \leq \frac{R^2}{\gamma^2}$$

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- 4: observe x_t, y_t
- 5: predict $\hat{y}_t = \arg \max_{y \in \mathcal{Y}(x)} w_t^\top \phi(x_t, y)$ \leftarrow this is expensive
- 6: **if** $\hat{y}_t \neq y_t$ **then**
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Let's look again at the structured SVM model from the previous lecture

StructSVM Objective

$$R(w) = \frac{1}{N} \sum_{i=1}^{N} \left(\max_{\mu} (w^{\top} F \cdot \mu + l_i^{\top} \mu) - w^{\top} F_i \cdot \mu_i \right) + \frac{\lambda}{2} ||w||^2$$

Rewriting in the same notation as perceptron:

StructSVM Objective

$$R(w) = \frac{1}{N} \sum_{i=1}^{N} \left(\max_{y \in \mathcal{Y}} (w^{\top} \phi(x_t, y) + \ell(y_t, y)) - w^{\top} \phi(x_t, y_t) \right) + \frac{\lambda}{2} ||w||^2$$

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- R(w) is convex (sum of affine & convex functions)
- denote $\tilde{y} = \arg \max_{y} (w^{\top} \phi(x_t, y) + \ell(y_t, y))$
- can be minimized using batch subgradient method

$$\frac{\partial R}{\partial w} = \frac{1}{N} \sum_{i=1}^{N} \left(\phi(x_t, y_t) - \phi(x_t, \hat{y}) \right) + \lambda w$$
$$w_{t+1} = w_t - \alpha_t \frac{\partial R}{\partial w}$$
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- R(w) is convex (sum of affine & convex functions)
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Where do we get \tilde{y} ?

$$\tilde{y}_i = \operatorname*{arg\,min}_{y \in \mathcal{Y}} (w^{\top} \phi(x_i, y) - \ell(x_i, y))$$

- this is the loss-augmented inference task
- requires a solution to a search problem in the underlying space
- has to be solved for every input instance (big data is a problem)
 - ⇒ solvable in $O(T \cdot K)$ on chains and trees (long sequences are a problem)
 - much harder or intractable for general graphs (e.g. MRFs)

So we have two problems:

- when there is relation to the inference, the learning loop is expensive
 - the non-probabilistic models above (Perceptron or SVM)
- or inference and learning maybe not related at all
 - ➡ e.g. CRF: the learning objective knows nothing about inference
 - same for many deep learning-based models
- often finding \tilde{y} alone is not enough
 - CRFs require feature expectations
 - large-margin methods require n-best lists
- Learning as Search Optimization (LaSo) is tackling these problems
 - get rid of the expensive $\arg \max$ (global decision)
 - decompose the structure building into local decisions

Setup:

- input $x \in \mathcal{X}$ induces a search space $\mathcal{Y}(x)$
- initial state b (also encodes x)
- set of states ${\cal S}$
- transition function $P(s_{t+1}|s_t, a_t)$ (here deterministic)
- for each (valid) sequence of states and actions, there is a corresponding output y(e)
- \blacksquare loss $\ell(e)=\ell(y^*,y(e)),$ where y^* is the ground truth structure
- feature generating function $\phi:\mathcal{S}\to\mathbb{R}^d,$ that expresses both the input x and previous actions
- agent follows a policy $\pi(a_t|s_t)$, which chooses an action a_t in state s_t
- trajectory $\tau_{\pi} = (a_1, s_1, \dots, s_{T-1}, a_T, s_T)$, where $P(s_{t+1}|s_t, \pi(s_t))$
- T is maximum lengths of au
- **goal**: $\min_{\pi} \mathbb{E}_{(x,y) \sim \mathcal{D}}[\ell(y, \tau_{\pi}))]$

Search space



[Daume'15]

- for any $s \in \mathcal{S}$ and $y \in \mathcal{Y}$ we can say if s can lead to y
- in this case we call it 'y-good', otherwise 'y-bad'
- ∎ goal:
 - ➡ first node can lead to any structure
 - ➡ the queue always contains at least one y-good node

Algo Learn(problem, initial, enqueue, w, x, y) $nodes \leftarrow MakeQueue(MakeNode(problem, initial))$ while *nodes* is not empty do $node \leftarrow \text{RemoveFront}(nodes)$ if none of $nodes \cup \{node\}$ is y-good or GoalTest(node) and node is not y-good then $sibs \leftarrow siblings(node, y)$ $\boldsymbol{w} \leftarrow update(\boldsymbol{w}, x, sibs, node \cup nodes)$ $nodes \leftarrow MakeQueue(sibs)$ else if GoalTest(node) then return w $next \leftarrow Operators(node)$ $nodes \leftarrow enqueue(problem, nodes, next, w)$ end if end while

- online learning algorithm
- search and learning are tightly intervened
- usually needs a size restriction on the queue
- similar to structured perceptron, except for early updates
- updates are done on errors (if the current beam cannot lead to the right answer)

$$w_{t+1} = w_t + \sum_{n \in sibs} \frac{\phi(x, n)}{|sibs|} - \sum_{n \in nodes} \frac{\phi(x, n)}{|nodes|}$$

Analysis:

- basically, follows the analysis for structured perceptron
- **•** result: for separable problems with margin γ , number of errors is $\leq \frac{R^2}{\gamma^2}$
- actually it's not exactly true for subtle reasons, see [Xu and Fern, 2007]

- natural only for sequence labelling problems
- hard to apply for tasks with production in arbitrary order

Literature



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