

Consistency and nonparametric algorithms

Consistency is an asymptotical property certifying that the risk of the predictors generated by a learning algorithm converges to the Bayes risk in expectation as the size of the training set increases. Recall that $A(S_m)$ is the predictor generated by a learning algorithm A on a training set S_m of size m . A learning algorithm A is **consistent** with respect to a loss function ℓ if for any data distribution \mathcal{D} it holds that

$$\lim_{m \rightarrow \infty} \mathbb{E}[\ell_{\mathcal{D}}(A(S_m))] = \ell_{\mathcal{D}}(f^*)$$

where the expectation is with respect to the random draw of the training set S_m of size m from the distribution \mathcal{D} , and $\ell_{\mathcal{D}}(f^*)$ is the Bayes risk for (\mathcal{D}, ℓ) . In some cases, we may define consistency with respect to a restricted class of distributions \mathcal{D} . For example, in binary classification we may restrict to all distributions \mathcal{D} such that $\eta(\mathbf{x}) = \mathbb{P}(Y = 1 \mid \mathbf{X} = \mathbf{x})$ is a Lipschitz function on \mathcal{X} . Formally, there exists $0 < c < \infty$ such that

$$|\eta(\mathbf{x}) - \eta(\mathbf{x}')| \leq c \|\mathbf{x} - \mathbf{x}'\| \quad \text{for all } \mathbf{x}, \mathbf{x}' \in \mathcal{X}.$$

Technically, this conditions implies that η is Lipschitz in \mathcal{X} . This is a restriction on the set of all allowed η as $c < \infty$ implies continuity (but the opposite is not true).

Nonparametric algorithms. Given a learning algorithm A , let \mathcal{H}_m be the set of predictors generated by A on training sets of size m : $h \in \mathcal{H}_m$ if and only if there exists a training set S_m of size m such that $A(S_m) = h$. We say that A is a nonparametric learning algorithm if A 's approximation error vanishes as m grows to infinity. Formally,

$$\lim_{m \rightarrow \infty} \min_{h \in \mathcal{H}_m} \ell_{\mathcal{D}}(h) = \ell_{\mathcal{D}}(f^*).$$

Two notable examples of nonparametric learning algorithms are k -NN and the greedy algorithm for decision tree classifiers (i.e., the algorithm that always chooses to split a leaf that maximizes the decrease in training error). Nonparametric algorithms are recognizable because:

- the size (memory footprint) of their predictors tends to grow with the training set size
- for any m and for all S_m , $\min_{h \in \mathcal{H}_m} \ell_S(s)$ is close to zero.

The standard k -NN algorithm is nonparametric but not known to be consistent for any fixed value of k . Indeed, one can only show that

$$\lim_{m \rightarrow \infty} \mathbb{E}[\ell_{\mathcal{D}}(k\text{-NN}(S_m))] \leq \ell_{\mathcal{D}}(f^*) + 2\sqrt{\frac{\ell_{\mathcal{D}}(f^*)}{k}} \quad (1)$$

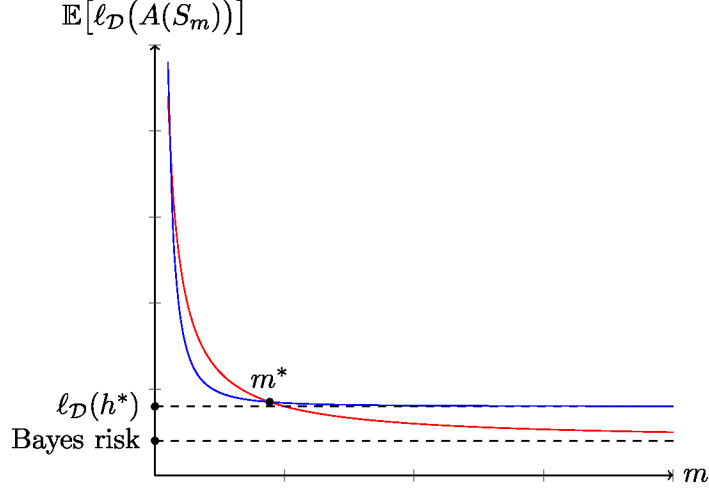


Figure 1: Typical behavior of expected risk $\mathbb{E}[\ell_{\mathcal{D}}(A(S_m))]$ as a function of training set size for a consistent algorithm (red line) and for a nonconsistent algorithm (blue line). For small training set sizes $m < m^*$, the nonconsistent algorithm has a better performance. (Thanks to Edoardo Marangoni for drawing the picture.)

for any data distribution \mathcal{D} . However, if we let k be chosen as a function k_m of the training set size, then the algorithm becomes consistent provided $k_m \rightarrow \infty$ and $k_m = o(m)$.

Similarly, the greedy algorithm for building tree classifiers is consistent (for $\mathcal{X} \equiv \mathbb{R}^d$) whenever the two following conditions are fulfilled. Let $\ell(\mathbf{x})$ be the leaf to which $\mathbf{x} \in \mathbb{R}^d$ is routed in the current tree and let N_ℓ be the number of training examples routed to a leaf ℓ . Then, as $m \rightarrow \infty$, to guarantee consistency we must have that

1. $\text{diameter}\left(\left\{\mathbf{x}' \in \mathbb{R}^d : \ell(\mathbf{X}) = \ell(\mathbf{x}')\right\}\right) \rightarrow 0$
2. $N_{\ell(\mathbf{X})} \rightarrow \infty$

where both conditions must hold when $m \rightarrow \infty$ and in probability with respect to the random draw of \mathbf{X} . In other words, the tree must grow unboundedly but not too fast.

In practice, a consistent algorithm may not be preferred over a nonconsistent one, see Figure 1. This due to the fact that the rate of convergence to the Bayes risk of a consistent algorithm can be arbitrarily slow, as shown by the following result.

Theorem 1 (No Free Lunch). *For any sequence a_1, a_2, \dots of positive numbers converging to zero and such that $\frac{1}{16} \geq a_1 \geq a_2 \geq \dots$ and for any consistent learning algorithm A for binary classification with zero-one loss, there exists a data distribution \mathcal{D} such that $\ell_{\mathcal{D}}(f^*) = 0$ and $\mathbb{E}[\ell_{\mathcal{D}}(A(S_m))] \geq a_m$ for all $m \geq 1$.*

Theorem 1 does not prevent a consistent algorithm from converging fast to the Bayes risk for

specific distributions \mathcal{D} . What the theorem shows is that if A converges to the Bayes risk for any data distribution, then it will converge arbitrarily slow for some of these distributions.

For binary classification, we can summarize the situation as follows.

- Under no assumption on η , there is no guaranteed convergence rate to Bayes risk.
- Under Lipschitz assumptions on η , the typical convergence rate to Bayes risk is $m^{-1/(d+1)}$.
- Under no assumptions on η , the typical convergence rate to the risk of the best predictor in a parametric (or finite) class \mathcal{H} is $m^{-1/2}$, exponentially better than the nonparametric rate.

Note that the convergence rate $m^{-1/(d+1)}$ implies that to get ε -close to Bayes risk, we need a training set size m of order $\varepsilon^{-(d+1)}$. This exponential dependence on the number of features of the training set size is known as **curse of dimensionality** and refers to the difficulty of learning in a nonparametric setting when datapoints live in a high-dimensional space.