



Practical ML Advice

Based on slides from Jude Shavlik and
Tom Dietterich

Proper Experimental Methodology Can Have a Huge Impact:

A 2002 paper in *Nature* (a major journal) needed to be corrected due to “training on the testing set”

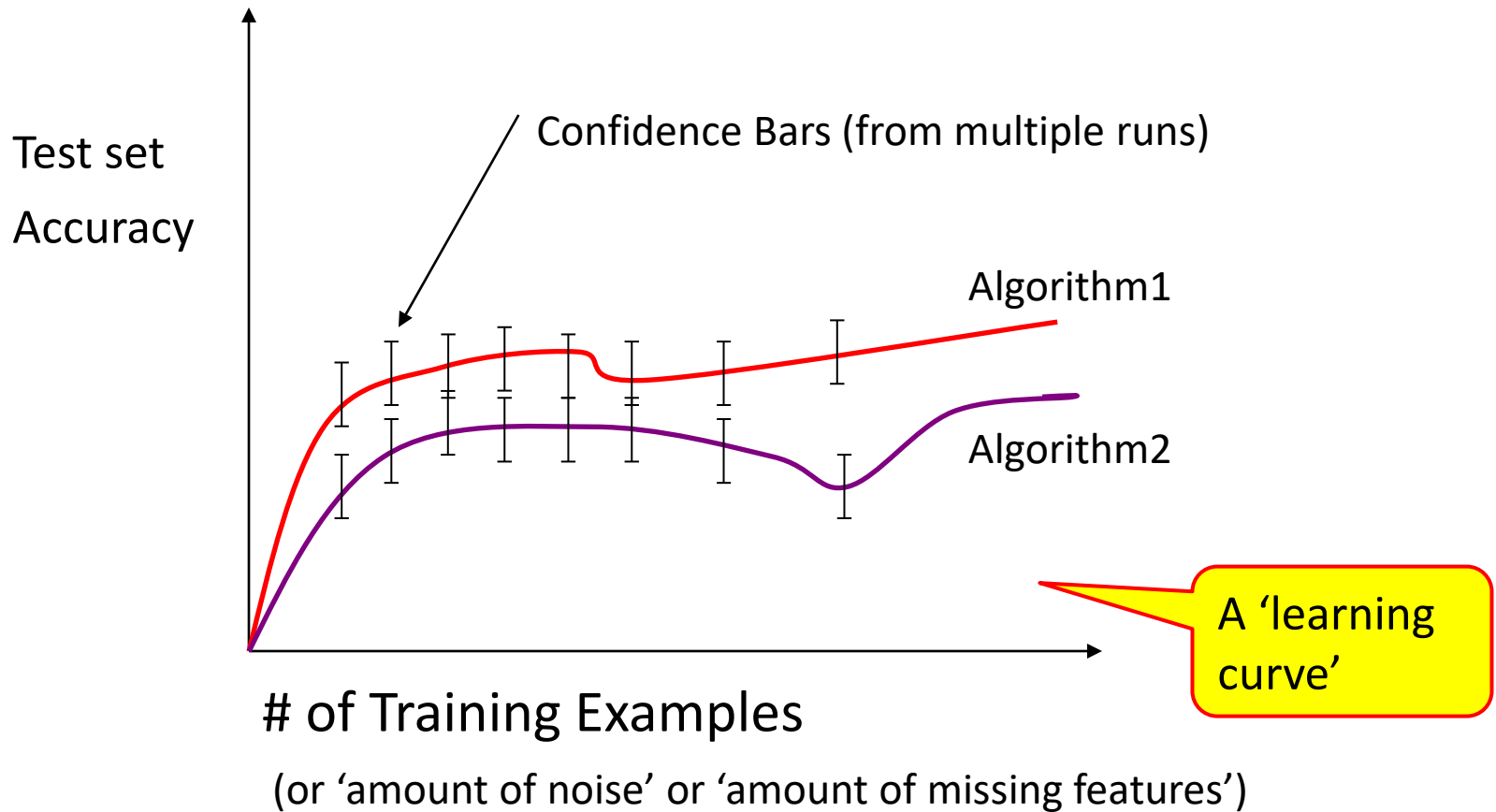
Original report : 95% accuracy (5% error rate)

Corrected report (which still is buggy):

73% accuracy (27% error rate)

Error rate increased over 400%!!!

Some Typical ML Experiments



Typical Experiments



	Test Set Performance
Full System	80%
Without Module A	75%
Without Module B	62%

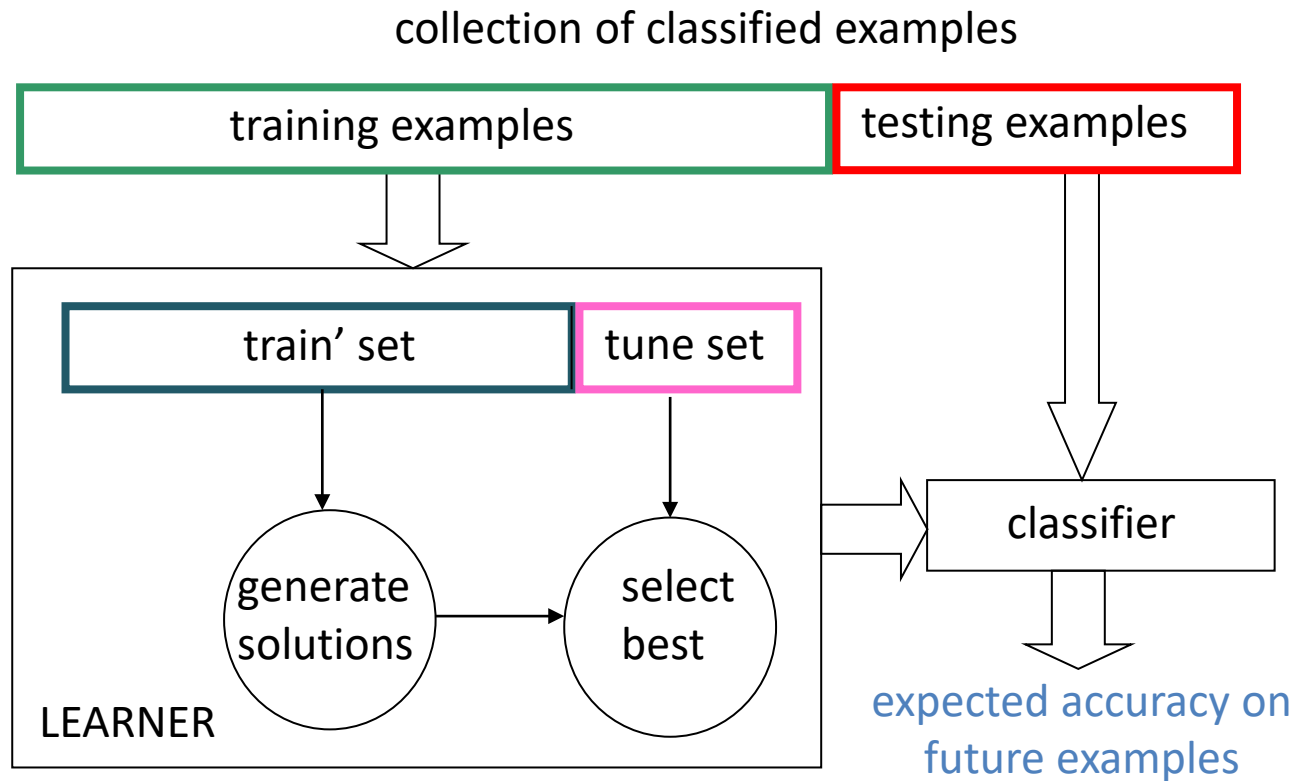
- 1) Start with a dataset of labeled examples
- 2) Randomly partition into N groups
- 3a) N times, combine $N - 1$ groups into a train set
- 3b) Provide **training set** to learning system
- 3c) Measure accuracy on “left out” group (the **test set**)



Called **N -fold cross validation**

- Often, an ML system has to choose when to stop learning, select among alternative answers, etc.
- One wants the model that produces the highest accuracy on **future** examples (“overfitting avoidance”)
- It is a **“cheat”** to look at the **test** set while still learning
- Better method
 - Set aside part of the training set
 - Measure performance on this validation data to estimate future performance for a given set of hyperparameters
 - Use best hyperparameter settings, train with **all** training data (except **test** set) to estimate future performance on **new** examples

A typical Learning system



Statistical techniques such as 10-fold cross validation and *t*-tests are used to get meaningful results

Multiple Tuning sets



- Using a **single** tuning set can be unreliable predictor, plus some data “wasted”
 - 1) For each possible set of hyperparameters
 - a) Divide training data into **train** and **valid.** sets, using ***N*-fold cross validation**
 - b) Score this set of hyperparameter values: average **valid.** set accuracy over the *N* folds
 - 2) Use **best** set of hyperparameter settings and **all** (train + valid.) examples
 - 3) Apply resulting model to **test** set



EVALUATING ML MODELS

Contingency Tables



(special case of 'confusion matrices')

		True Answer	
		+	-
Algorithm Answer	+	$n(1,1)$ [true pos]	$n(1,0)$ [false pos]
	-	$n(0,1)$ [false neg]	$n(0,0)$ [true neg]

Counts of occurrences

A 2x2 contingency table with 'Algorithm Answer' on the y-axis and 'True Answer' on the x-axis. The y-axis has labels '+' and '-'. The x-axis has labels '+' and '-'. The four cells contain: top-left: n(1,1) [true pos]; top-right: n(1,0) [false pos]; bottom-left: n(0,1) [false neg]; bottom-right: n(0,0) [true neg]. Two blue arrows originate from the bottom center of the table, pointing to the bottom-left and bottom-right cells.

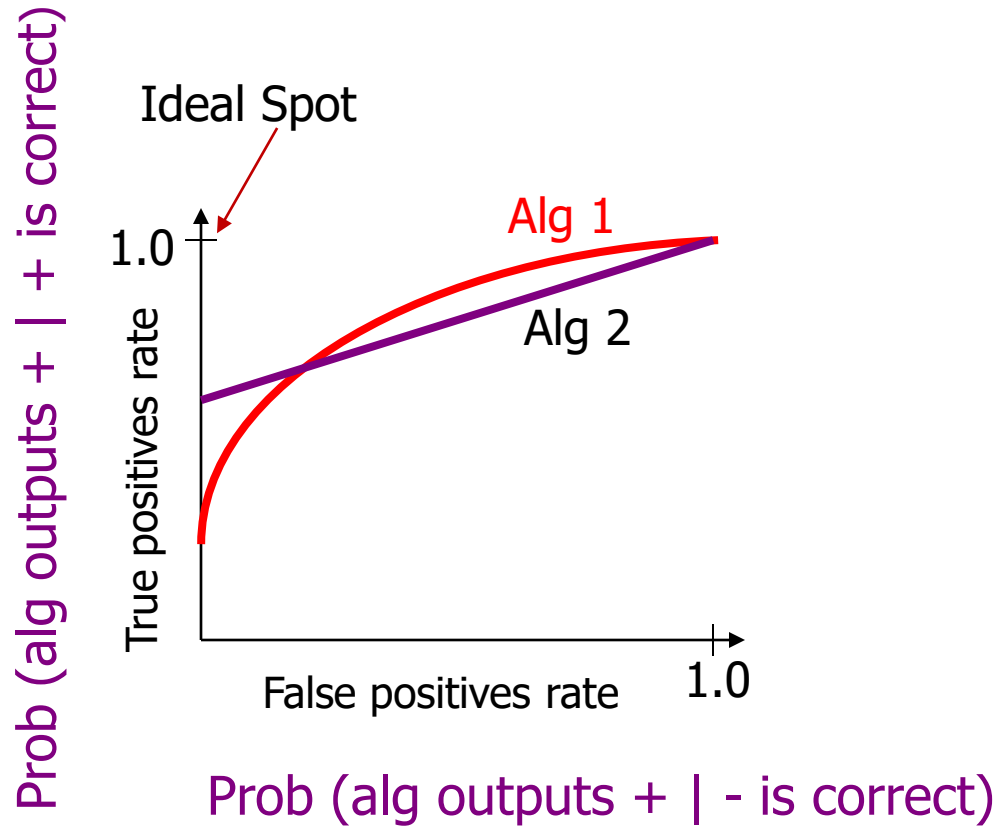
True Positive Rate
(TPR) = $n(1,1) / (n(1,1) + n(0,1))$
= correctly categorized +'s / total positives
~ $P(\text{algo outputs } + \mid + \text{ is correct})$

False Positive Rate
(FPR) = $n(1,0) / (n(1,0) + n(0,0))$
= incorrectly categorized -'s / total neg's
~ $P(\text{algo outputs } + \mid - \text{ is correct})$

Can similarly define False Negative Rate and True Negative Rate

- ROC: *Receiver Operating Characteristics*
- Started for radar research during WWII
- Judging algorithms on accuracy alone may not be good enough when **getting a positive wrong** costs more than **getting a negative wrong** (or vice versa)
 - e.g., medical tests for serious diseases
 - e.g., a movie-recommender system

ROC Curves Graphically



Different algorithms can work better in different parts of ROC space. This depends on cost of false + vs false -

Creating an ROC Curve



The Standard Approach:

- You need an ML algorithm that outputs **NUMERIC** results such as prob(example is +)
- You can use ensemble methods to get this from a model that only provides Boolean outputs
 - e.g., have 100 models vote & count votes

Alg. for Creating ROC Curves

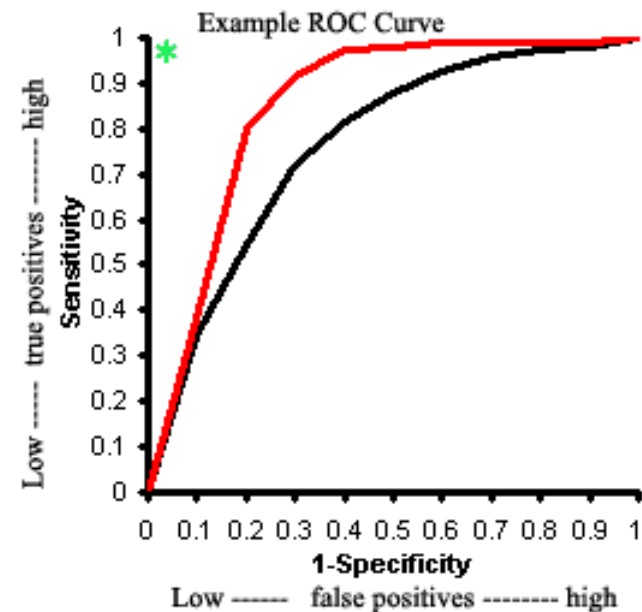


Step 1: Sort predictions on test set

Step 2: Locate a *threshold* between examples with opposite categories

Step 3: Compute TPR & FPR for each threshold of Step 2

Step 4: Connect the dots

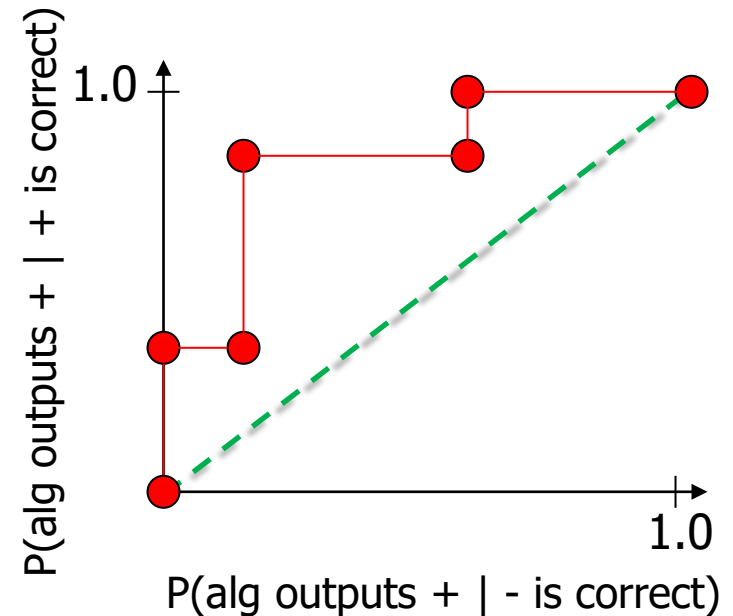


Plotting ROC Curves - Example



ML Algo Output (Sorted) Correct Category

Ex 9	.99		+
Ex 7	.98	TPR=(2/5), FPR=(0/5)	+
Ex 1	.72	TPR=(2/5), FPR=(1/5)	-
Ex 2	.70		+
Ex 6	.65	TPR=(4/5), FPR=(1/5)	+
Ex 10	.51		-
Ex 3	.39	TPR=(4/5), FPR=(3/5)	-
Ex 5	.24	TPR=(5/5), FPR=(3/5)	+
Ex 4	.11		-
Ex 8	.01	TPR=(5/5), FPR=(5/5)	-

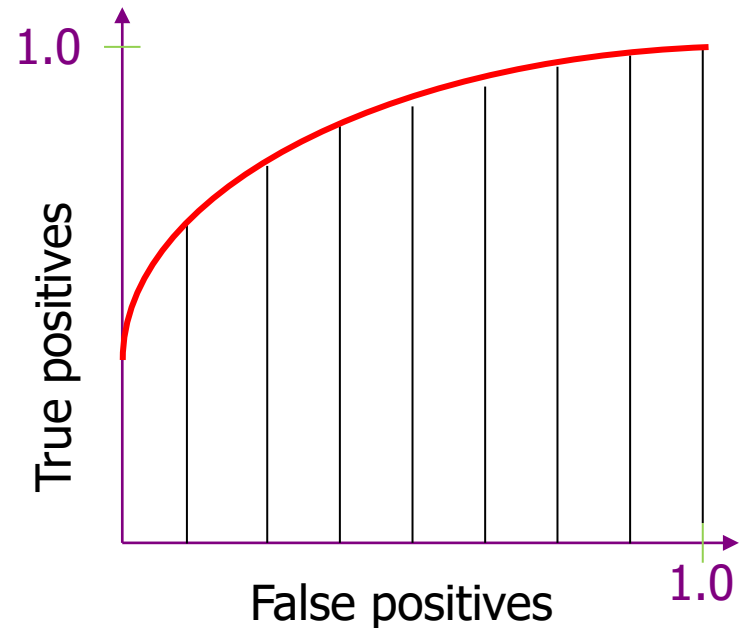


Algorithm predicts + if its output is ≥ 0

Area Under ROC Curve



- A common metric for experiments is to numerically integrate the ROC Curve
 - Usually called AUC
 - Probability that ML alg. will “rank” a randomly chosen positive instance higher than a randomly chosen negative one
 - Given a randomly selected positive example and a randomly selected negative example, AUC is the probability that the classifier will be able to distinguish them
- Can summarize the curve **too much** in practice



- One strength of ROC curves is that they are a good way to deal with **skewed** data ($|+| \gg |-|$) since the axes are fractions (rates) independent of the # of examples
- You must be careful though!
 - Low FPR * (many negative ex) = **sizable number of FP**
 - Possibly more than # of TP

Precision vs. Recall

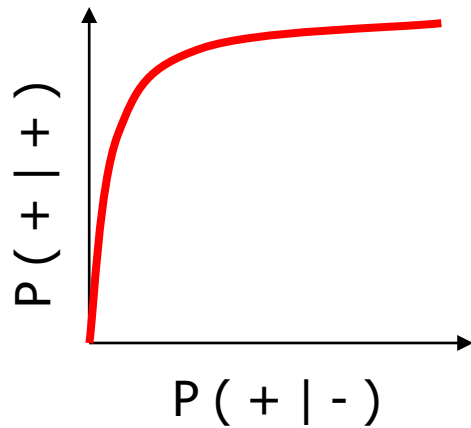


- Think about search engines...
- **Precision** = (# of relevant items retrieved)
/ (total # of items retrieved)
= $n(1,1) / (n(1,1) + n(1,0))$
- **Recall** = (# of relevant items retrieved)
/ (# of relevant items that exist)
= $n(1,1) / (n(1,1) + n(0,1))$
= TPR
- Notice that $n(0,0)$ is not used in either formula
Therefore you get no credit for filtering out irrelevant items

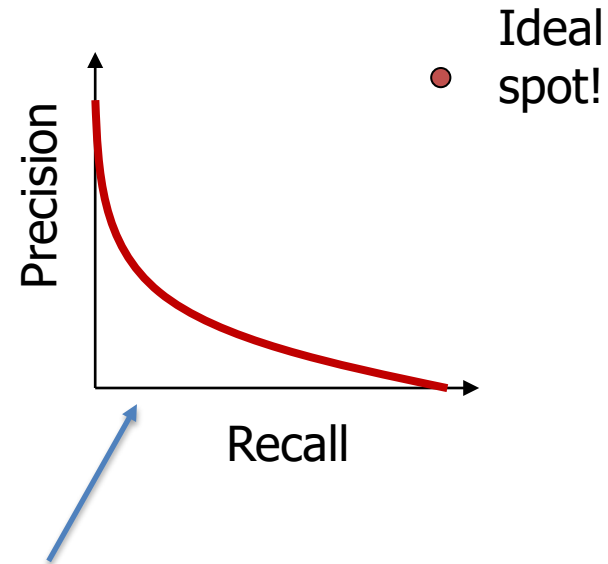
ROC vs. Precision-Recall



You can get very different visual results on the same data!



vs.

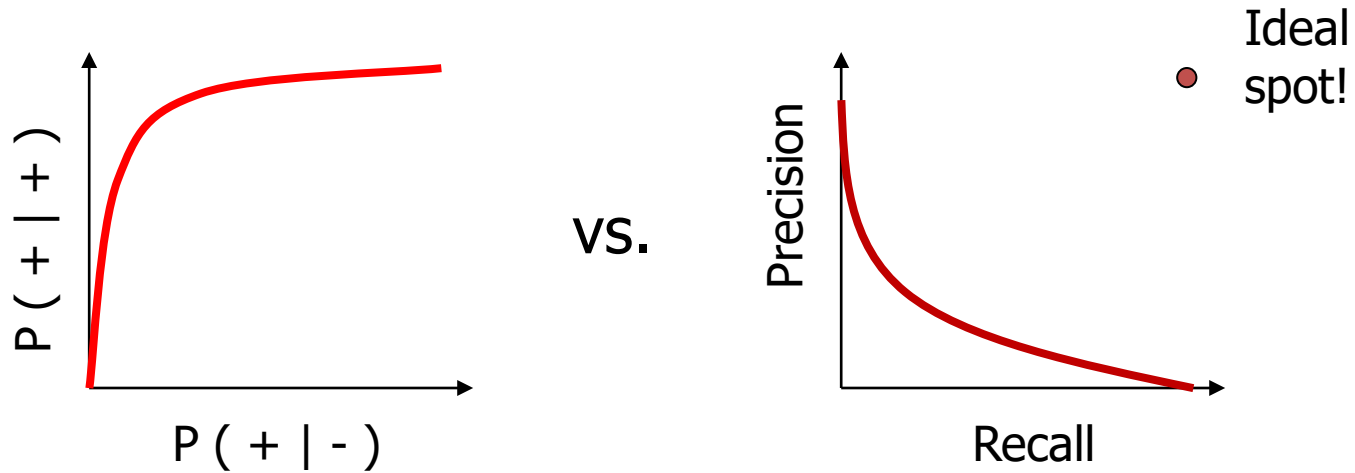


Produced by varying threshold for positive identification, e.g., say 1 if $p(1|x) > .5$ in logistic regression

ROC vs. Precision-Recall



You can get very different visual results on the same data!



The reason for this is that there may be lots of $-$ ex's (e.g., might need to include 100 neg's to get 1 more pos)

Rejection Curves

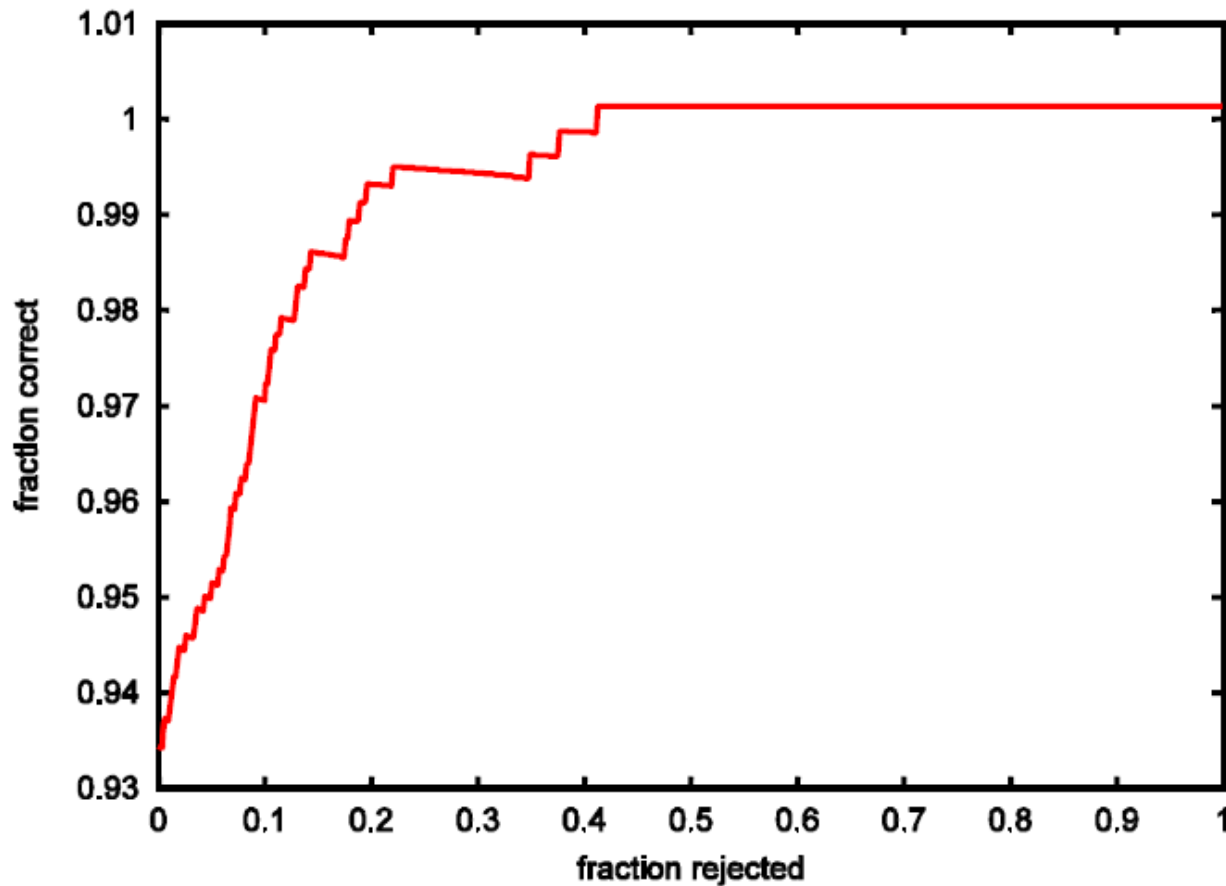


- In most learning algorithms, we can specify a threshold for making a rejection decision
 - Probabilistic classifiers: adjust cost of rejecting versus cost of FP and FN
 - Decision-boundary method: if a test point x is within θ of the decision boundary, then reject
 - Equivalent to requiring that the “activation” of the best class is larger than the second-best class by at least θ

Rejection Curves



- Vary θ and plot fraction correct versus fraction rejected



The F1 Measure



- Figure of merit that combines precision and recall

$$F_1 = 2 \cdot \frac{P \cdot R}{P + R}$$

where P = precision; R = recall. This is twice the harmonic mean of P and R .

- We can plot F_1 as a function of the classification threshold θ