Rule-based Regression

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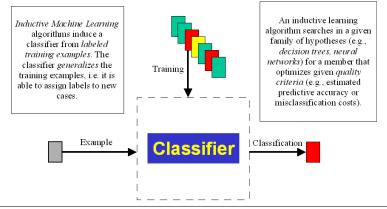
Machine Learning



- Definition (Mitchell, 1997)
 - "A computer program is said to *learn* from experience E with respect to some class of tasks T and performance measure P, if its performance at tasks in T, as measured by P, improves with experience E."
- Given:
 - ► a task T
 - ► a performance measure P
 - some experience E with the task
- Goal:
 - generalize the experience in a way that allows to improve your performance on the task

Indroduction of Classifiers





Introduction of Classifiers



The most "popular" learning problem:

- ► Task:
 - learn a <u>model</u> that predicts the outcome of a dependent variable for a given instance
- Experience:
 - experience is given in the form of a data base of examples
 - an example describes a single previous observation
 - instance: a set of measurements that characterize a situation
 - Iabel: the outcome that was observed in this situation
- Performance Measure:
 - compare the predicted outcome to the observed outcome
 - estimate the probability of predicting the right outcome in a new situation

Data Representation

Attribute-Value Data



- Each example is decribed with values for a fixed number of <u>attributes</u> (also called <u>features</u>)
 - Nominal Attributes:
 - store an unordered list of symbols (e.g., color)
 - Numeric Attributes:
 - store a number (e.g., income)

A sample task



Day	Temperature	Outlook	Humidity	Windy	Play Golf?
07-05	26	sunny	high	false	no
07-06	28	sunny	high	true	no
07-07	29	overcast	high	false	yes
07-09	23	rain	normal	false	yes
07-10	20	overcast	normal	true	yes
07-12	12	sunny	high	false	no
07-14	8	sunny	normal	false	yes
07-15	25	rain	normal	false	yes
07-20	18	sunny	normal	true	yes
07-21	18	overcast	high	true	yes
07-22	20	overcast	normal	false	yes
07-23	19	rain	high	true	no
07-26	11	rain	normal	true	no
07-30	16	rain	high	false	yes

today	9	sunny	normal	false	?
tomorrow	13	sunny	normal	false	?

A sample task



Day	Temperature	Outlook	Humidity	Windy	Play Golf?
07-05	26	sunny	high	false	no
07-06	28	sunny	high	true	no
07-07	29	overcast	high	false	yes
07-09	23	rain	normal	false	yes
07-10	20	overcast	normal	true	yes
07-12	12	sunny	high	false	no
07-14	8	sunny	normal	false	yes
07-15	25	rain	normal	false	yes
07-20	18	sunny	normal	true	yes
07-21	18	overcast	high	true	yes
07-22	20	overcast	normal	false	yes
07-23	19	rain	high	true	no
07-26	11	rain	normal	true	no
07-30	16	rain	high	false	yes

possible rules:

play=no \leftarrow temperature ≥ 25.5 \land temperature < 28.5 $play=no \leftarrow temperature < 14$ \wedge temperature \geq 9.5 $play=no \leftarrow outlook=rainy \land$ windy=true

today	9	sunny	normal	false	?
tomorrow	13	sunny	normal	false	?

A sample task



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07-23	19	rain	high	true	no
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possible rules:

play=no \leftarrow temperature > 25.5 \land temperature < 28.5 $play=no \leftarrow temperature < 14$ \land temperature > 9.5 $play=no \leftarrow outlook=rainy \land$ windy=true

but also (t=temperature):

 $\begin{array}{l} \mathsf{play=no} \leftarrow t < 26.5 \land t \geq \\ 25.5 \land \mathsf{outlook=sunny} \land \\ \mathsf{humidity=high} \land \mathsf{windy=false} \\ \mathsf{play=no} \leftarrow t < 28.5 \land t \geq \\ 27.5 \land \mathsf{outlook=sunny} \land \\ \mathsf{humidity=high} \land \mathsf{windy=true} \\ \cdots \end{array}$

today	9	sunny	normal	false	?
tomorrow	13	sunny	normal	false	?

Separate-and-conquer Rule Learning



- Separate-and-conquer (or Covering) paradigma (originated from the AQ algorithm (Michalski, 1969))
- ▶ still used in most Rule Learning systems (e.g., RIPPER (Cohen, 1995))
- 1. Generalization: extend the current theory by a "good" rule
- 2. Separate: remove all examples covered by this rule
- 3. Conquer: if examples left, goto 1.
- rules are combined in a decision list
 - sorted list of rules
 - the first rule that "covers" the example is used to classify the example
 - if no rule covers the example the last rule is used as a default rule (predicts the majority class)

Searching for a single rule



- generate the first rule that covers all examples
- generate all refinements of the current rule by creating all attribute-value pairs from the data
 - nominal attributes: use equality tests (i.e., =)
 - \blacktriangleright numerical attributes: use inequality tests (i.e., \geq and <)
- add each refinement to the current rule and test which is the best for a given (heuristic) criterion
- if a new best is found store it
- if the error of the rule is 0 stop the process and return the best rule that was found during this process

Combining rules in a decision list



- if a rule is found add the rule to the sorted list of rules
- remove all the examples that are covered by the rule
- ▶ if all but the remaining *n* examples are covered stop inducing rules (currently n = 1)
- <u>else</u>: search for the next rule on the remaining examples
- as last rule add a default rule that predicts the majority class

Rule Learning Heuristics



- Rule Learning Heuristics implement the criterion for evaluating rules
- many Rule Learning Heuristics for classification are known (based on positive and negative examples)
- Parametrized trade-off between
 - Consistency: (1 error) of the rule and
 - Coverage: how many examples are covered by the rule
- Heuristics for Regression (positive and negative examples are not known here) rely on
 - the current error/loss (Consistency in classification) of the rule
 - the coverage of the rule
- Regression Heuristics may also feature a parameter that trades off between the error and the Coverage of the rule

From Classification to Regression



- instead of predicting a discrete outcome in Regression the outcome is continuous
- 2 ways to deal with this:
 - 1. discretize numeric outcome and use standard classification algorithms
 - problem: number of classes has to be known in advance
 - algorithm used to discretize: P-CLASS (Weiss and Indurkhya, 1995))
 - 2. adapt the algorithm to Regression tasks
 - example for an adaption in Rule Learning
 - either predict a certain value (*Median* or *Mean*) in the head of the rule directly (like we did)
 - or use a (linear) model in the head to predict the value (algorithm M5RULES (Holmes, Hall, and Frank, 1999), (Quinlan, 1992))

Regression measures



- Deviation from Mean $def = \frac{1}{n} \sum_{i=1} (y_i y')^2$
- Normalized Mean Squared Error NMSE = MSE/def
- Relative Coverage RC = COVERAGE(r)/n
- ► Relative Cost Measure $h_{rcm} = c \cdot (1 NMSE) + (1 c) \cdot RC$

where n = # of examples left, y_i = true value, \bar{y}_i = predicted value, y' = mean of all instances, r = the current rule

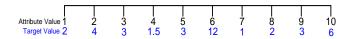
Current implementation



- numerical and nominal attributes, numerical target variable
- covering paradigma
- interchangeable heuristics and splitpoint computing methods
- parameters:
 - parameter of the heuristic
 - parameter for splitpoint computation
 - to reduce the number of splitpoints for a numerical attribute a clustering was used
 - the parameter determines how many clusters are computed
 - percentage of coverage of ruleset (for inducing the default rule)
 - currently all but the last remaining example has to be covered

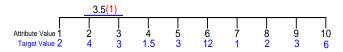


- if all possible splitpoints (those between 2 instances) for all numeric attributes are used the search space explodes
- remedy: do not create all splitpoints but cluster examples together that minimize some error criterion
- ▶ and use only the splitpoints between these clusters (currently about 5-10)
- Algorithm:
 - sort the examples of the attribute in ascending order
 - remove duplicates by setting the mean over all duplicates as target value
 - merge examples that minimize the mean absolute error



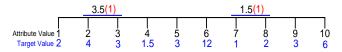


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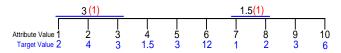


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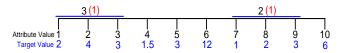


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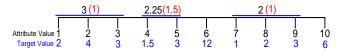


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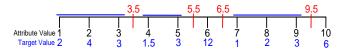


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Evaluation of the models



- ▶ for domain-dependent evaluation we used *MAE* and *RMSE* = \sqrt{MSE}
- for domain-independent evaluation we used the correlation coefficient (between predicted and actual value)
- we also record model complexity by measuring the number of rules and conditions (for rule based models)
- 1x10 cross-validation with same folds for each model
- ▶ our approach was compared to M5RULES, LINEAR REGRESSION, SVMREG (all implemented in *weka* (Witten and Frank, 2005))

Results

In terms of MAE



- preliminary results (sp = 10, c = 0.45) for 13 datasets from the UCI-Repository (Asuncion and Newman, 2007)
- second number describes standard deviation among the 10 folds of the CV

dataset	SeCo	M5Rules	Linear Regression	SVMReg
auto-horse	16.61 ± 6.35	15.85 ± 10.25	13.64 ± 3.24	13.48 ± 4.0
auto-mpg	4.44 ± 1.49	3.03 ± 0.81	2.87 ± 0.98	2.83 ± 0.98
auto-price	2526.6 ± 773.1	2157.8 ± 937.4	2450.5 ± 1084.0	2292.32 ± 1012.05
breast-tumor	8.02 ± 0.73	7.79 ± 0.74	7.9 ± 0.72	8.2 ± 0.76
cloud	0.45 ± 0.15	0.3 ± 0.12	0.27 ± 0.07	0.28 ± 0.09
сри	36.80 ± 29.38	15.19 ± 9.17	47.7 ± 20.89	24.95 ± 23.52
echo-month	13.41 ± 2.62	8.68 ± 2.97	8.48 ± 3.15	9.08 ± 2.73
housing	5.43 ± 2.98	3.39 ± 1.44	3.99 ± 2.13	3.73 ± 2.05
meta	95.59 ± 170.29	232.52 ± 190.24	146.54 ± 148.08	96.91 ± 166.08
sensory	0.64 ± 0.13	0.73 ± 0.14	0.76 ± 0.18	0.77 ± 0.19
servo	0.54 ± 0.15	0.32 ± 0.11	0.62 ± 0.12	0.53 ± 0.17
strike	274.61 ± 116.46	287.0 ± 87.31	264.73 ± 84.0	228.49 ± 83.23
veteran	91.28 ± 59.05	92.99 ± 44.4	92.99 ± 44.4	82.58 ± 54.89
average rank	3.08	2.27	2.58	2.08

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Results

In terms of different parametrizations



- the number of splitpoints are fixed to 10 but the parameter of the heuristic is varied
- Iowest errors are marked blue

dataset	c = 0.45		<i>c</i> = 0.5		<i>c</i> = 0.6		c = 0.7	
	MAE	# rules	MAE	# rules	MAE	# rules	MAE	# rules
auto-horse	16.6	2	15.2	16	21.4	35	16.5	57
auto-mpg	4.44	1	3.92	157	3.62	184	3.64	226
auto-price	2526	6	2922	7	3104	46	2836	48
breast-tumor	8.0	0	8.5	13	10.7	209	10.4	236
cloud	0.45	7	0.46	6	0.42	12	0.39	42
сри	36.8	5	37.8	7	38.8	9	29.3	15
echo-month	13.4	0	14.2	79	14.2	92	13.2	87
housing	5.43	5	4.7	43	4.54	369	4.47	427
meta	95.6	3	95.2	30	147.8	69	147	124
sensory	0.63	0	0.82	430	0.86	404	0.9	428
servo	0.54	4	0.39	20	0.39	22	0.38	29
strike	274	0	362	234	361	300	368	359
veteran	91	0	116	70	119	82	123	91

Discussion



- our algorithm implements a Separate-and-conquer Regression Rule Learner
- trade-off between consistency and coverage is more complex than it is in classification
 - tuning of the parameters has to be analyzed better
- but the current implementation is competetive to other rule-based implementations (that do not predict models in the head)
- a new splitpoint computing method was introduced
 - only about 10 splitpoints are sufficient for most of the datasets
 - much more faster than computing all splitpoints
 - but optimal cluster number still has to be found

Future Work



> this is work-in-progress so there are many ways to improve the algorithm

- by determine a suitable setting of the cluster parameter
- by systematically tune the parameter of the heuristic
 - previously we tuned the parameters of 5 heuristics for classification
 - we also want to find the best parameter for regression
- by avoiding overfitting by leaving more examples uncovered
- predict (linear) models in the head of the rule
- try to visualize the behaviour of the different heuristics in a space similar to Coverage Spaces
- include domain-independent comparison with RRMSE = $\sqrt{\frac{MSE}{def}}$

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