

Rule Induction Algorithm for Application to Geological and Petrophysical Data

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Abstract: Very large geological, geophysical, and petrophysical databases often contain multiple data types that must be interpreted for application to subsurface modeling. Significant advances in discovering complex and even nonintuitive data relationships could lead to better predictions. There is a litany of data analysis techniques used today, including cluster analysis, principal component analysis, discriminant analysis, parametric and nonparametric regression, and N -dimensional histograms. Regression techniques and neural networks have in common their multivariate combination of predictor variables. These techniques may be good at interpolating within the data boundaries of the training data, but may be poor for extrapolation because of the lack of understanding of the underlying relationships in the variables. Alternatively, machine learning and data mining technologies including Rough Sets hold the promise of finding data category relationships and expressing those in a rule-based language. This paper presents a novel rule induction algorithm derived from these machine-learning techniques, developed for reservoir characterization with geological and geophysical data. A set of facies models with systematical changing in the geometric features is synthesized. The geometric features are coded and the effective permeability is calculated. Rules between effective permeability and geometric features are deduced by using the proposed technique. The consistence of the deduced rules with those implemented in the data synthesization exhibit the effectivity of the proposed technique. Further a second example of facies assignment from wireline logs is used to test the proposed technique. The deduced rules are confirmed by the geologists who spend significant time trying to summarize rules from the well logs. The probability feature of the rules and the distinguishability analysis feature of the proposed technique supplied additional information for the geologist to reconsider their original distinction among facies.

1 Background

The field of data mining [2, 7, 8, 11, 17, 18] has grown in recent years to deal with large databases available in different industries, in particular, the financial and medical fields. Data mining is the identification or discovery of patterns in data. There are several different types of data mining. These include classification,

clustering (segmentation), association, and sequence discovery. The main focus of classification is supervised induction, that is, inference of rules and relationships from large databases. The aim is to extract knowledge from data, so that results not directly in the training data set can be predicted. The training data helps to distinguish predefined classes. Neural networks [3, 9, 13, 14, 27], decision trees [5, 6, 16, 19, 26, 29] and if-then-else rules are classification techniques. A disadvantage of neural networks is that it is difficult to provide a good rationale for the predictions made, that is, the rules are not always clear.

Data mining is an interdisciplinary field bringing together techniques from statistics, machine learning, artificial intelligence, pattern recognition, database, and visualization technologies. The methods used in data mining are not fundamentally different from older quantitative model-building techniques, but are natural extensions and generalizations of such methods especially applicable to uncontrolled, big real world data. Like in other fields, petroleum industry involves extra large data collection routinely and the data collection is not for satisfying the requirement of data analysis purpose. The subsurface feature and the complexity of petroleum reservoir make data collected noisy, incomplete, and redundant, and there are valuable information remaining undiscovered in the data. The general goal of data mining is to extract implicit, previously unknown, hidden, and potentially useful information from raw data. In many cases traditional techniques may not be efficient or not suitable. For example, strong statistical assumptions and the adherence to a specific statistical distribution model may make statistical methods not suitable. The inability to generalize knowledge, usually represented in the form of rules, of traditional techniques may not fulfill the requirement from industries to derive meaningful relationship of variables from raw data. Data mining as a distinct discipline to attack the challenges of deriving knowledge from large database has potential for the knowledge discovery in the petroleum reservoir characterization. There are applications of various data mining techniques to petroleum characterization [1, 4, 12, 15, 28].

A rule-based algorithm is intended to provide understandable rule-like relationships in the data. A rule is a prevailing quality or state. Induction is an instance of reasoning from a part to a whole. Rules indicate the degree of association between variables, map data into predefined classes, and identify a finite set of categories or clusters to describe the data. The rules support specific tasks and are generated by repeated application of a certain technique, or more generally an algorithm, on the data. Rough Sets [17, 20-25, 30, 31] are specialized methods for inducing rules. The essential idea of rough sets is to express uncertain knowledge through an approximation space, which is constructed as certain sets.

Many other methods, including regression analysis, assume that there is a functional form between the predictor and response variables. These smooth out variations and are difficult to apply to multivariate nonlinear responses. Discriminant analysis separates samples into groups based on relationships in the

training data. The relations must be linear combinations of variables that are made explicit. N -dimensional histograms are used to delineate a relationship between a response and multiple predictors, which preserves the uncertainty in the relation by reading the value of the response directly from the set of predictors. Principal components analysis is a popular technique to discover the source of variation within the data but again results are expressed as linear combinations of multiple variables. All of these techniques can “fit” the data but might not be good predictors and do not provide insight into relationships. Clustering is a method to partition the database into segments where each segment member shares similar qualities. Clustering techniques may include optimization algorithms to determine the maximum similarity among members within each group and a minimum similarity among members across the groups.

Obvious applications of rule induction in petroleum geological data mining include (1) predicting reservoir facies from multiple wireline logs, (2) predicting reservoir permeability from wireline logs, (3) predicting reservoir facies from multiple seismic attributes, and (4) predicting stratigraphic geometries and spatial architecture from quantified analog, outcrop, seismic, and numerical stratigraphic data.

2 Rule Induction Algorithm

In rough set theory, concepts of *coverage* and *accuracy* are used to describe uncertainty in probabilistic rules. In practice, every object in the data table may lead to a rule. The challenge is to identify significant rules, that is, accurate rules with high frequency of occurrence. Data mining has no *a priori* model, that is, no functional relationship between the data is assumed.

A feature of geological data is that most rules are probabilistic and not deterministic. Another aspect of geological data is that there are relatively few attributes from a classical data mining perspective. These two considerations were used to develop the concepts and terms described below. A complete definition of the terms is in the glossary.

2.1 Data Table

Table 1 is a schematic *data table* for rule induction. The table has M observations and N condition attributes. Without loss of generality, it is assumed that there is a single decision attribute. Multiple decision attributes merely increase the number of rules and the configuration space. Initially, the condition attributes and decision attribute can have different values or categories, which must be transformed into discrete codes. The codes are based on statistical analysis of the data for each attribute. For example, the continuous variable of “gamma-ray count” may be transformed into 3 discrete codes of *high*, *average*, and *low*.

Table 1. A typical data table for rule induction.

	ca_1	ca_2	...	ca_{N-1}	ca_N	de
o_1	0	2	...	1	2	1
o_2	2	1	...	1	2	1
...
o_{M-1}	2	2	...	0	1	0
o_M	1	0	...	2	2	2

ca , condition attribute; de , decision attribute; o , observation;
 N , number of condition attributes; M , number of observations

The number of categorical values that condition attribute i , ca_i , can take is n_i , ($i = 1, \dots, N$), n_o is the number of categorical value the decision attribute de can take. The number of discrete or categorical values the decision attribute and some condition attributes can take is fixed, or will be determined for a problem, say $n_1 = 3$, but the codes themselves can be different; both $\{0,1,2\}$ and $\{5,9,11\}$ are valid codes for the three categorical values of attribute 1.

2.2 Configuration and Data Coverges

A *configuration* is a unique combination of the values of condition attributes. The total number of possible configurations is: $N_N = \prod_{i=1}^N n_i$, where n_i , ($i = 1, \dots, N$) is the number of categorical values of condition attribute i . Note that some configurations may not appear in the training data table. Further, define $N_i \left(= \prod_{k=1}^i n_k \right)$ as the cumulative number of configuration up to condition attribute i , $i = 1, \dots, N$. Configuration index j , $j = 1, \dots, N_N$, is uniquely determined by the values l_i , $i = 1, \dots, N$ taken by each condition attribute, where $j = \sum_{i=1}^N (l_i - 1) \times \frac{N_N}{N_i} + 1$, and $l_i \in 1, \dots, n_i$. The condition attribute values l_i , $i = 1, \dots, N$ from a configuration index j may be retrieved easily.

Every configuration may lead to potential rules in the system. Since the decision attribute can take n_o outcomes, there are $N_N \times n_o$ potential rules for the system.

The occurrence of configuration j with outcome o is counted as $C_{j,o}$, (j , $j = 1, \dots, N_N$ and o , $o = 1, \dots, n_o$) from the data table. If $C_{j,o} = 0$, then there are no observations in the data table corresponding to that configuration and outcome pair. The number of observations M is:

$$M = \sum_{j=1}^{N_N} \sum_{o=1}^{n_o} C_{j,o}$$

The *configuration coverage* C_j is defined as the total number of observations associated with configuration j , *i.e.*,

$$C_j = \sum_{o=1}^{n_o} C_{j,o}, \quad j = 1, \dots, N_N$$

The *decision coverage* D_o is specified as the number of observations associated with outcome o , *i.e.*,

$$D_o = \sum_{j=1}^{N_N} C_{j,o}, \quad o = 1, \dots, n_o$$

The *relative (configuration) coverage* of configuration j , \hat{C}_j is a measure used to determine closeness to sufficiency of observations in obtaining a reliable rule. \hat{C}_j is undefined for $C_j = 0$, equal to zero for $C_j = 1$, and asymptotic to one as the *coverage* increases:

$$\hat{C}_j = 1 - \frac{1}{C_j}, \quad j = 1, \dots, N_N$$

The definition of *relative (configuration) coverage* could be different.

The *relative (decision) coverage* of a decision value, $\hat{D}_{j,o}$, is a measure of the proportion of the occurrence frequency of a configuration and decision outcome pair out of the number of observations associated with outcome o , *i.e.*,

$$\hat{D}_{j,o} = \frac{C_{j,o}}{D_o}, \quad j = 1, \dots, N_N, o = 1, \dots, n_o$$

These relative coverages will be used to adjust the importance of rules. For example, more weight will be placed on decision values with small global proportions to avoid cases where $C_{j,o}$ is small, but is close to D_o . In such a case, this configuration is very important to that decision.

2.3 Conditional Probability and Accuracy

The *conditional probability* of each outcome value, o , ($o = 1, \dots, n_o$), for configuration j , ($j = 1, \dots, N_N$) is defined as:

$$p_{o|j} = \frac{C_{j,o}}{C_j}$$

Since there are n_o possible decision attribute categories, a probability of $1/n_o$ implies no information. Any conditional probability different from $1/n_o$ entails preference in the decision category, *i.e.*, some knowledge. Specifically, as $p_{o|j}$ nears 1, configuration j implies decision category o . A closeness of $p_{o|j}$ to 0 means configuration j does not lead to decision category o . Conditional probabilities close to 1 or 0 contain equally important information for rule induction. The former leads to a positive rule relating configuration j to a specific decision category o and the latter leads to negative rules relating configuration j to some other decision category. For a data system, a N_N by n_o conditional probability table will be established.

The knowledge conveyed by the conditional probability leads us to define a measure of *accuracy* as follows:

$$a_{o|j} = \begin{cases} p_{o|j} \cdot n_o - 1, & \text{if } p_{o|j} \leq \frac{1}{n_o} \\ \frac{(p_{o|j} \cdot n_o - 1)}{n_o - 1}, & \text{otherwise} \end{cases}$$

The accuracy $a_{o|j} = 0$ if $p_{o|j} = 1/n_o$ (no knowledge); the accuracy $a_{o|j} = 1$ when we have $p_{o|j}$ equals to 1 (perfect positive rule); and, $a_{o|j} = -1$ when we have $p_{o|j}$ equals to 0 (perfect negative rule). A graph of accuracy versus probability is shown on Figure 1.

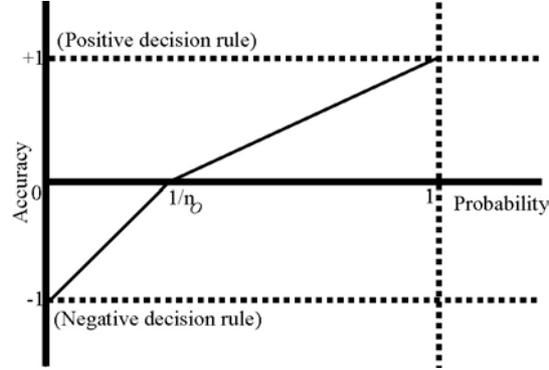


Fig. 1. Relationship between accuracy and conditional probability.

2.4 Significance

The *significance* is defined as a combined measure of the *accuracy* and *coverage*, which is used to provide a relative ranking to the rules. The significance is the product of the *accuracy* and the *relative configuration coverage*:

$$S_{j,o} = \hat{C}_j \cdot a_{o|j}$$

This equation holds for negative accuracy. For positive accuracy, the significance is further modified to account for the *relative decision coverage*:

$$S_{j,o} = \hat{C}_j \cdot a_{o|j} \cdot \hat{D}_{j,o}$$

As mentioned above, $\hat{D}_{j,o}$, is used to account the proportion of occurrence of an outcome with a specific configuration to the overall occurrence of that outcome, which serves like a normalization factor considering the unequally occurrence of decision attribute categories. Positive accuracy results from occurrence exceeding the average occurrence and the scaling using $\hat{D}_{j,o}$ will bring up the importance for rules associated with low occurrence outcome categories. Negative accuracy always results from occurrence lower than average level, the scaling factor is not applied to avoid bringing the significance towards zero. For example, a zero occurrence $C_{j,o}$ leads to negative one in accuracy (perfect negative rule) and large negative significance value, applying $\hat{D}_{j,o}$ will revert it to zero significance.

2.5 Rule Table

Rules are the configurations and decision outcome pairs with large positive or negative significance. Positive *rules* are extracted from the configuration and decision outcome pairs ranked from positive 1 to zero by positive *significance*, and negative rules are extracted from the configuration and decision outcome pairs ranked from negative one to zero *significance*. A practitioner can decide whether to assign a threshold for significance. Positive rules provide evidence for the decision outcome and negative rules provide evidence against the decision. If a configuration and decision outcome pair's significance does not meet the criteria, the rule table entry is "missing" for lack of data.

We expect many "missing" entries for rule induction with real data. The "missing" entries will be eliminated or reduced based on compatible rules generated from subsets of data variables.

2.6 Distinguishability

In a data mining exercise, there is an implicit assumption that all decision outcomes are distinguishable from the data. This assumption is not always true and a measure of *distinguishability* can assist in determining the decision that is distinct on the basis of the data table.

Table 2. Predicted probability from the rules.

Observation (m)	Predicted probability from the	real outcome (\tilde{o}_m)
1	$p_{1.1}, p_{1.2}, \dots, p_{1.n}$	\tilde{o}_1
2	$p_{2.1}, p_{2.2}, \dots, p_{2.n.}$	\tilde{o}_2
...
m	$p_{m.1}, p_{m.2}, \dots, p_{m.n.}$	\tilde{o}_{M-1}
...
$M-1$	$p_{M-1.1}, p_{M-1.2}, \dots, p_{M-1.n.}$	\tilde{o}_{M-1}
M	$p_{M.1}, p_{M.2}, \dots, p_{M.n.}$	\tilde{o}_M

Once the rule induction is finished, one can construct a data table like the one shown in Table 2. For each observation in the training data set, a predicted probability is assigned by looking for the right configuration j and reading the corresponding conditional probability p_{oj} .

The data table may be divided into n_o classes with an indicator function:

$$ind(m; o) = \begin{cases} 1 & \text{if } o = \tilde{o}_m, \quad m = 1, \dots, M; \quad o = 1, \dots, n_o \\ 0 & \text{otherwise} \end{cases}$$

where \tilde{o}_m is the true outcome category for observation m .

The summation of $ind(m; o)$ is a counting of all outcomes related to decision value o :

$$D_o = \sum_{m=1}^M ind(m; o)$$

The average predicated probability associated with prediction of o is:

$$\hat{p}_o = \frac{1}{D_o} \sum_{m=1}^M ind(m; o) \cdot p_{m,o}, \quad o = 1, \dots, n_o$$

where $p_{m,o}$ is the predicted probability value of observation m having decision category o .

As \hat{p}_o approaches 1, prediction is good. As \hat{p}_o goes to $p_o = \frac{D_o}{M}$, *i.e.*, the global proportion for decision outcome o , there is no information provided by the rules.

Therefore, the relative information value for prediction outcome o is defined as:

$$I_o = \frac{\hat{p}_o - p_o}{p_o}$$

A composite information value can be defined as the average over all outcomes:

$$\bar{I} = \frac{1}{n_o} \sum_{o=1}^{n_o} I_o$$

The expected predicated probability and information measure is used to determine whether outcomes o and o' are distinct.

Let's consider merging two different decision outcomes o and o' , then we have $n_o - 1$ new decision outcomes and the predicted probability table (like Table 2) will be updated. Updating the predicted probability is straightforward. We derive

$n_o - 1$ probability values from the previous n_o values. One of them will be $p_{o|j} + p_{o'|j}$ when the actual outcome turns out to be o or o' .

Following the same procedure, new I_o and the \bar{I} values can be calculated. If we define $\bar{I}_{o:o'}$ as the information value with o and o' merged, the change in \bar{I} is:

$$\Delta\bar{I}_{o:o'} = \bar{I}_{o:o'} - \bar{I}$$

$\Delta\bar{I}_{o:o'} > 0$ means that o and o' merged leads to an improved prediction. Thus, one should consider whether to treat outcomes o and o' together, *i.e.*, as indistinguishable, since the given training data table cannot differentiate them. One could seek to obtain additional data on a different data type that would distinguish those outcomes. $\Delta\bar{I}_{o:o'} < 0$ means that merging the outcomes leads to a poorer prediction. Thus, the outcomes should be treated as distinct.

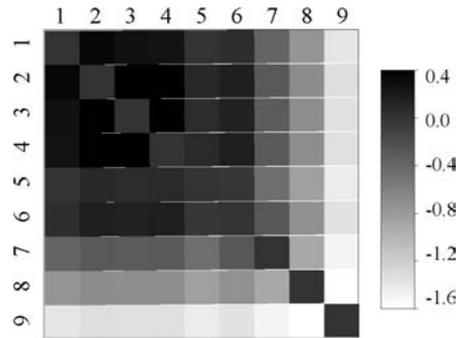


Fig. 2. Information changes for lumped/merged decision categories.

Lumping or *merging* could be considered for all possible pairs of outcome values, and the $\Delta\bar{I}$ can be tabulated, plotted or ranked from high to low. The high positive values are candidates for merging while the low positive and negative values should be kept separate. The $\Delta\bar{I}_{o:o'}$ table can be visualized as a grayscale map. For example, Figure 2 shows a map of the matrix pairing nine decision outcomes from the facies assignment example. The bigger the $\Delta\bar{I}$ value, the higher the distinguishability of outcome pair and the darker the color. The diagonal cells are the pairing of outcomes with themselves and have $\Delta\bar{I}$ values of

0. As shown in the Figure, outcomes 7, 8, 9 are very distinguishable from other outcomes. The upper left corner area with positive $\Delta\bar{I}$ values indicate that outcomes 1-2, 2-3, and 2-4 are not as distinguishable and might be candidates for combination.

2.7 Data or Condition Attribute Value

The same measure of information can be used to consider the value (importance) of condition attributes or sets of condition attributes. All procedures described above are based on the entire data set with all condition attributes, but the procedures can be applied to any data subset with partial condition attributes.

For a N condition attribute data set, there are $2^N - 1 (= \sum_{l=1}^N C_N^l)$, (l is the number of condition attributes in the subset and $l = 1, \dots, N$) subsets. For each subset, a new data set is derived and the algorithm is applied. A new set of configurations of condition attributes is obtained and p_{oj} values are calculated for every decision category o for a given configuration j . For each individual decision category, the expected predication probability of the occurrence of decision category o when given the data entry having decision category o is also calculated as well as the information measure. The information value for each subset is, for example, $\bar{I}_{\{1,2,3,\dots,N\}} = \bar{I}$, for all attributes, $\bar{I}_{\{2,3,\dots,N\}}$ for leaving out condition attribute 1, $\bar{I}_{\{2\}}$ for just attribute 2, *etc.* These results can be plotted as a graph or tabulated in a table that ranks the value of the condition attributes (subsets). The largest values, or maximums of information in the subsets with the same number of condition attributes, would normally tend to increase as the number of attributes in the subsets increases. However, noisy data or attributes that do not yield additional information value could reduce \bar{I} .

2.8 Completion of Rule Table

The rule (conditional probability) tables built from the entire data set are not complete since there may be no observations for some condition attribute configurations and there are some rules have been screened out due to small significance. A complete rule table is necessary for practical usage of rule induction algorithm and “missing” entries in the rule table should be filled in. Filling “missing” entries in the rule table will be accomplished based on significant rule derived from data subsets.

There are many subsets with various number of condition attributes, therefore, it is necessary to order the subsets in filling procedure. This ordering will be based on their information value. The ordered subsets will be considered until all “missing”

entries are filled in. The entries filled from subsets are flagged to indicate the subset used.

The utilization of the rules is two-fold: (1) gain insight and make general observations from the rule table, and (2) use the conditional probability associated with each configuration and decision outcome pair in the conditional probability table to model the predictions probabilistically. These applications will be illustrated with two examples.

The probabilities are used in modeling. Of course, there is opportunity to reduce uncertainty by incorporating additional data sources in the modeling phase.

3. Implementation

There are issues related to (1) the physical meaning and subjective understanding of the rules, (2) the usage of results, (3) the meaning of a probabilistic prediction, and (4) data integration. The significance has value between -1 to + 1. In principle, the potential rules with largest absolute significance values will be chosen as important rules, but there exists ambiguity of how to define big and small, *e.g.*, is 0.9 close enough? When there are -1 and/or +1 in the significance list, those rules will be chosen without question. The others will depend on judgment. Fortunately, we are not concerned with rules in the middle of the list.

The rule induction is only as good as the training data. But in contrast to other techniques, *e.g.* neural networks, each rule is clear and understandable to a practitioner as to its applicability. Another issue is the selection of categorical classes by application of thresholds to a continuous variable. This could be performed by optimization.

The rule induction procedure consists of the following steps: (1) prepare the data table and determine optimal classes for the conditioning attributes (discretization), (2) enumerate all possible configurations and outcomes, (3) calculate the *conditional probability*, *configuration coverage*, *decision coverage*, *accuracy*, and *significance*, (4) present a table for each outcome $o = 1, \dots, n_o$ with the configurations sorted from low to high *significance*, (5) apply cutoff criteria on significance to build a rule table, (6) evaluate information value for the condition attributes and/or decision outcome, and (7) fill up the “missing” entries in the rule table based on subsets of the conditioning attributes.

4 Examples

4.1 Synthetic Fluvial Channel System

In fluvial or deepwater depositional setting, the sandy facies occur as sand channels with associated levee and crevasse deposits. For simplicity without loss generality, we only consider sand-filled channels embedded with a matrix of shale in this example. The corresponding effective permeability depends on the orientation, sinuosity, width, thickness and geometry of the channels. The relationship between these geometric features and the effective permeability are *rules* to be extracted for future use. For this purpose, channel facies models are created with a variety of channel parameters such as *orientation*, *thickness*, and *sinuosity*, which are called condition attributes for the rule induction system. Flow simulation is conducted for each generated model and the effective permeability in X direction is taken as quality measure (decision categories) of the model.

The synthesized data set consists of the geometrical features of channel sands and the associated effective permeability. An object modeling program, *fluvsim* in *GSLIB* [10], was used to generate multiple facies models. The simulation domain is a 100 by 100 two-dimensional area. Three triangular distributions with non-overlapped value ranges are set for each of the three parameters, *i.e.*, orientation (O), sinuosity (S), and width (W) (width/thickness ratio). Totally, there are 27 combinations (sets) of the parameter distributions. For each set of parameter distributions, parameters were randomly drawn and 50 facies models were generated. Thus, there are a total of 1350(=50×27) different facies models created. For each facies model, permeability values of 100 and 1 were assigned to channel sand and shale, respectively and a flow simulation was conducted with no-flow boundaries, using the *GSLIB* [10] program *flowsim*. The effective permeability in the X direction was taken as a measure of the quality or productivity of each facies model.

The data generated were organized as a data table with 1350 rows, each representing one facies model, and 4 columns, where the first three denotes the values of three condition attributes, *i.e.*, the *orientation* (O), *sinuosity* (S), and *width* (W) (*width/thickness ratio*) of the channels, and the fourth column was the value of the decision attribute, *i.e.*, the *effective permeability* (K) in X direction.

Figure 3 shows one of 50 facies models for two of 27 parameter distributions. Figure 4 shows the composite histogram of effective permeability of the 1350 facies models and the histograms of effective permeability of 50 facies models from the two parameter distributions shown in Figure 3, respectively.

The rule induction method requires that the data be *binned* into categorical variables. It is reasonable and straightforward to classify the three geometrical parameters (condition attributes) into three categories consistent with the three

distributions. The three categorical values of *orientation* (\mathbf{O}) are defined as 2 (good: azimuth angle of 0 ± 10 , which is aligned with the X direction), 1 (medium: azimuth angle of 45 ± 10) and 0 (bad: azimuth angle of 90 ± 10 , which is perpendicular to the X direction). The three categorical values for *sinuosity* (\mathbf{S}) are denoted as 2 (high: high deviation 30 ± 5 and short length of 20 ± 5 units), 1 (medium: medium deviation 20 ± 5 and medium length of 35 ± 5 units), 0 (low: low deviation 10 ± 5 and long length of 50 ± 5 units). The three categorical values for *width* (\mathbf{W}) are assigned as 2 (large: 25 ± 5), 1 (medium: 15 ± 5) and 0 (small: $5 - 10$). Columns 2 to 4 of Table 3 list the codes of the parameters used in the generation of facies models.

Table 3. Parameter used in facies models generation and the statistic of the effective permeability values.

No ^a	\mathbf{O}	\mathbf{S}	\mathbf{W}^b	mean K	Std K	K ^{low}	K ^{med}	K ^{high} ^c
1	0	0	0	3.90	1.53	32 (64)	18 (36)	0 (0)
2	0	0	1	2.22	0.48	49 (98)	1 (2)	0 (0)
3	0	0	2	1.91	0.31	50(100)	0 (0)	0 (0)
4	0	1	0	8.50	2.57	4 (8)	46 (92)	0 (0)
5	0	1	1	5.14	2.53	23 (46)	26 (52)	1 (2)
6	0	1	2	4.29	2.87	34 (68)	16 (32)	0 (0)
7	0	2	0	9.40	2.30	1 (2)	49 (98)	0 (0)
8	0	2	1	9.69	4.26	3 (6)	39 (78)	8 (16)
9	0	2	2	8.16	4.75	14 (28)	30 (60)	6 (12)
10	1	0	0	8.47	2.75	2 (4)	47 (94)	1 (2)
11	1	0	1	6.94	5.05	23 (46)	21 (42)	6 (12)
12	1	0	2	4.55	4.01	34 (68)	12 (24)	4 (8)
13	1	1	0	8.21	1.95	0 (0)	50 (100)	0 (0)
14	1	1	1	7.88	3.28	6 (12)	42 (84)	2 (4)
15	1	1	2	5.28	3.49	25 (50)	23 (46)	2 (4)
16	1	2	0	5.02	1.00	5 (10)	45 (90)	0 (0)
17	1	2	1	5.66	1.76	6 (12)	44 (88)	0 (0)
18	1	2	2	5.97	2.52	15 (30)	35 (70)	0 (0)
19	2	0	0	17.46	2.17	0 (0)	2 (4)	48 (96)
20	2	0	1	18.74	4.05	0 (0)	9 (18)	41 (82)
21	2	0	2	21.19	5.80	0 (0)	6 (12)	44 (88)
22	2	1	0	8.66	1.44	0 (0)	50 (100)	0 (0)
23	2	1	1	8.54	2.64	3 (6)	46 (92)	1 (2)
24	2	1	2	9.08	4.17	8 (16)	34 (68)	8 (16)
25	2	2	0	3.97	0.77	29 (58)	21 (42)	0 (0)
26	2	2	1	3.66	1.06	35 (70)	15 (30)	0 (0)
27	2	2	2	3.23	0.85	42 (84)	8 (16)	0 (0)

a: index of configuration; b: \mathbf{O} , Orientation; \mathbf{S} , Sinuosity; \mathbf{W} , width; c: K^{Low}, k≤4; K^{Medium}, 4<k<14; K^{High}, k≥14.

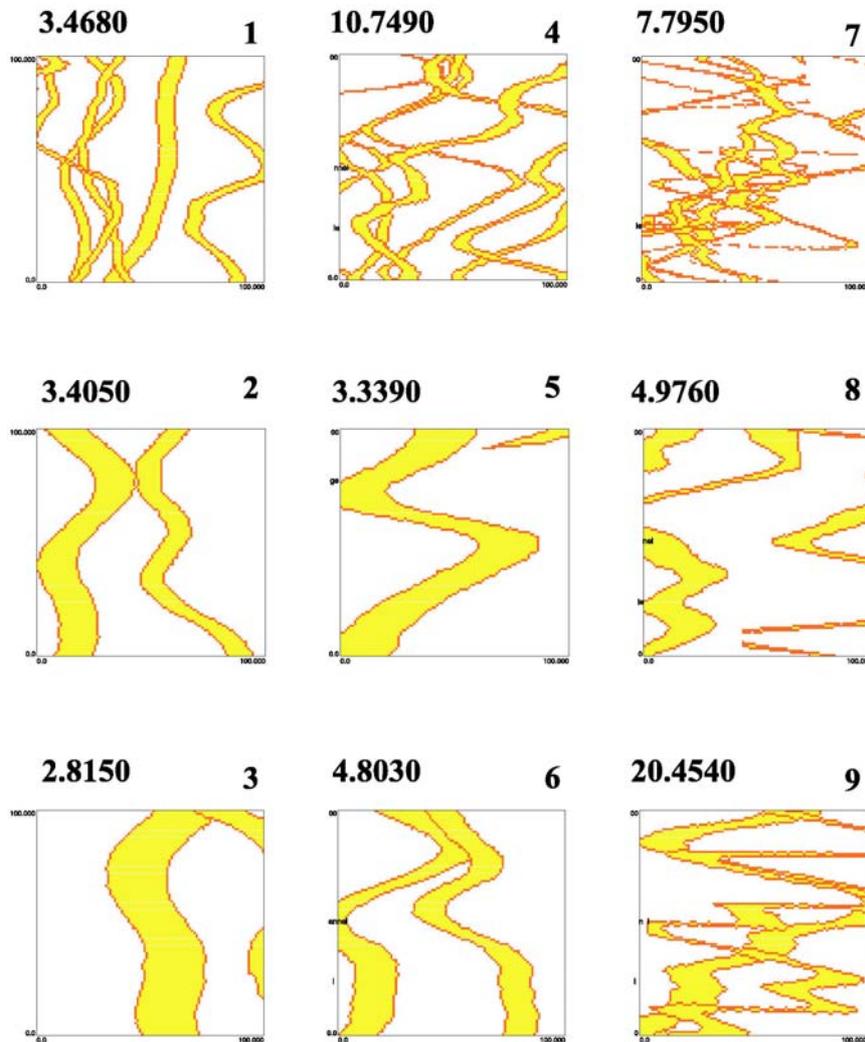


Fig. 3. Example geological models.

The continuous variable of effective permeability has to be discretized into categorical values. The method of *binning* the permeability data could affect the final rule induction results. For simplicity, three categories were determined for the effective permeability by just inspecting the permeability histograms. Summary statistics of the effective permeability for the 50 realizations of each configuration are listed in columns 5 and 6 of Table 3.

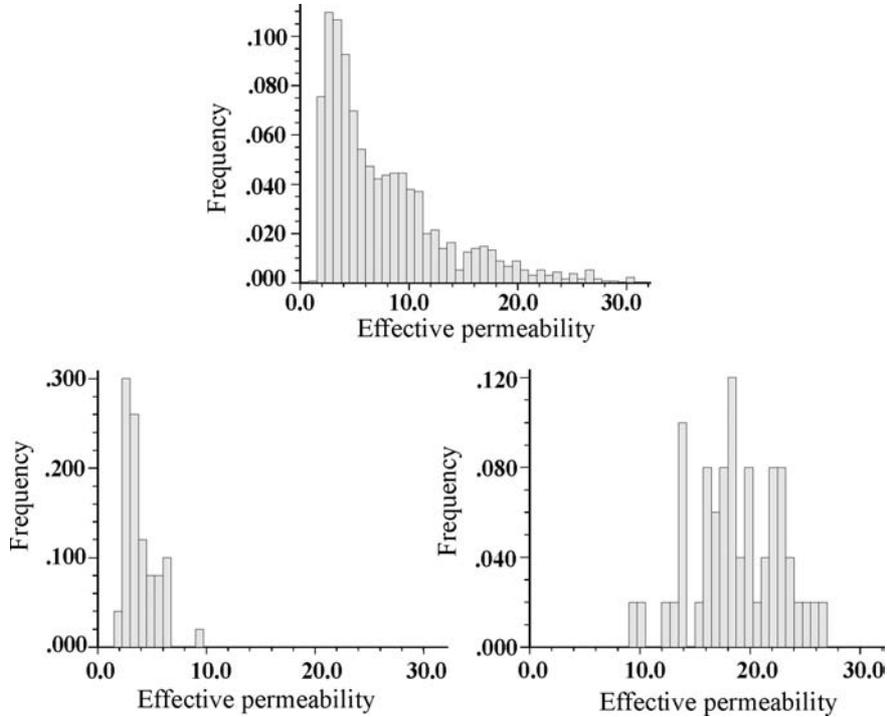


Fig. 4. *Left*, Composite histogram of effective permeability of overall 1350 facies model; histograms of effective permeability of 50 realizations for parameter distribution 1 (*middle*) and 20 (*right*).

Table 4. The expected rule for the fluvial system.

Good \mathcal{O} (2)	and	low \mathcal{S} (0)	and	No matter \mathcal{W}	leads to	high \mathcal{K} (2)
Bad \mathcal{O} (0)	and	low \mathcal{S} (0)	and	No matter \mathcal{W}	leads to	low \mathcal{K} (0)
Good \mathcal{O} (2)	and	High \mathcal{S} (2)	and	No matter \mathcal{W}	leads to	low \mathcal{K} (0)
Medium \mathcal{O}					mostly leads to	medium \mathcal{K}

The three categories determined for effective permeability are: 2 (high: $k \geq 14$), 1 (medium: $4 < k < 14$), and 0 (low: $k \leq 4$). The number of instances and proportions of the observations on each of the three decision categories for each configuration are tabulated in the last three columns of Table 3. By inspecting Table 3, we can expect certain rules like those listed in Table 4. Note that the “No Matter \mathcal{W} ” in Table 4 indicates that the width can have any of the categorical values.

4.2 Facies Assignment from Wireline Well Logs

This example consists of well logs data and facies assignment from cored wells. The purpose of the study is to investigate the potential to use well logs for facies assignment in uncored wells.

Table 5. Description of facies.

Facies	Description	Number of Data
1	Cross-bedded medium to coarse-grained sandstone	158
2	Medium to fine-grained sandstone	283
3	Well sorted fine-grained sandstone; shoreface	241
4	Very fine to fine-grained sandstone	181
5	Heterolithic: coarse-/fine-grained intervals; sand-filled burrow	117
6	Transition zone, lower shoreface to inner shelf	77
7	Muddy v. Fine-grained sandstone and sandy shale	62
8	Very fine sandy shale	70
9	Shale	27

Various well logs data are available including *Gamma Ray (GR)*, *Resistivity (Res)*, *Bulk Density (BD)* and *Neutron Activation (NA)*. Both *Bulk Density (BD)* and *Neutron Activation (NA)* are measures of density and the cross over of the difference between *normalized Bulk Density (BD)* and *normalized Neutron Activation (NA)* is important in well log interpretation in industry practice. Therefore, instead of using *Neutron Activation (NA)*, the difference of *normalized Bulk Density (BD)* and *Neutron Activation (NA)* are calculated as a new attribute, *CROSSOVER (CR)*. In addition, we have *water/oil/gas zone (Zone)* information for the wells and we use this piece of information as well. There are 9 facies recognized from cores and Table 5 lists the description of facies. The depth resolution in the wells is 0.5 feet. After removing missing data, there are a total of 1216 data points (observations) and the number of observations for each facies (decision value) is listed in Table 5 as well. Figure 5 shows the distribution of the four well log attributes of the entire data set.

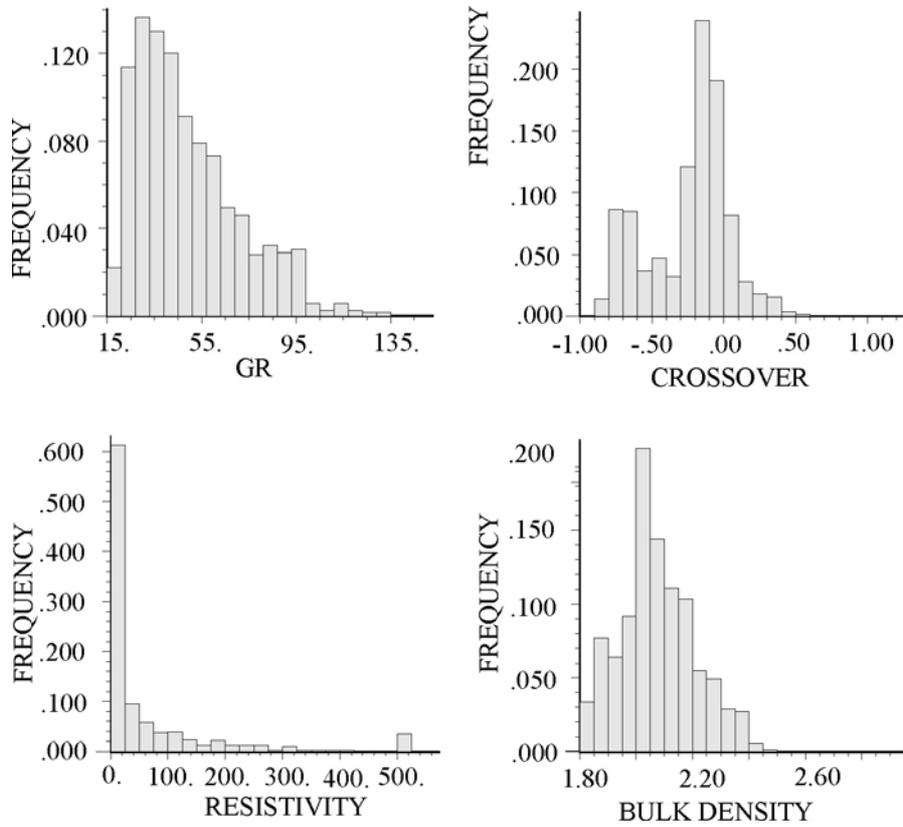


Fig. 5. Histograms of condition attributes.

Table 6. Codes of condition and decision attributes.

Code	0	1	2	3	4	5	6	7	8
GR	0-30	30-55	≥55						
Res	0-1.5	1.5-2.5	2.5-20	≥20					
BD	0-2.05	2.05-2.30	≥2.30						
CR	+	-0.5-0	≤-0.5						
Zone	0 (water)	1 (Oil)	2 (Gas)						
Facies	1	2	3	4	5	6	7	8	9

GR, Gamma ray; **Res**, Resistivity; **BD**, Bulk Density; **CR**, CrossOver; **Zone**, Water/Gas/Oil Zone

In order to use the rule induction algorithm, the four continuous attributes need to be discretized. This is not a trivial issue and research is ongoing for optimal discretization. We inspected and compared the distributions of the four attributes for each individual facies (decision attribute) to decide the discretization. The

discretization of the attributes and the coding are listed in Table 6. Note that we code all the attributes from 0. Finally a 1216 by 5 data table is constructed for rule induction.

5 Results, Conclusion and Discussion

5.1 Effective Permeability of Fluvial Sandy Channels

The algorithm was first applied to the 1350 by 4 synthesized channel sand/effective permeability data table and 27 configurations of conditioning attributes were extracted. Table 7 lists the configurations of condition attributes, the occurrence frequency $C_{j,o}$, and the significance $S_{j,o}$ of decision category o when given configuration j . By construction, all 27 possible configurations of condition attributes in this example have equal, non-zero sample coverage, *i.e.*, $C_j = 50$. This will not necessarily be the case for most data systems in practice.

The full set of rules is sorted by significance for each decision category, which is listed in Table 8. The positive rules shown in Table 9 are taken from the top portion of Table 8, which has the highest positive significance values. The negative rules as tabulated in Table 10 are taken from the bottom portion of Table 8, which have the lowest significance values.

The rule set is used to estimate the permeability class of training objects and Table 11 lists the accurate rate of the classification. Even though this accuracy rate derived from the training set is not a good measure to evaluate the predictability of the modeling, it still can provide some indication of the prediction model. The rules are quite reliable and predictable.

For this synthesized data set, the training data covers all configurations by construction. Therefore, the rule set is complete from the set with all condition attributes and there are no blanks in the rule sets to be filled. Figure 6 shows the information value of all 7 subsets of this synthesized data set and the full set with all three condition attributes has the highest information value. In the situation of the full set does not cover all configurations of condition attributes, such information values will be used to rank the subsets for retrieving corresponding rules from the subsets. Figure 7 shows the information value change when decision classes are lumped/merged pair-wise for the full condition attribute set. The diagonal elements set the basis for comparison that corresponds to a situation without decision-class lumping. Off-diagonal elements show the change in the information values when the attributes are lumped. Darker color than the diagonal elements indicate an increase in the value of information and those lighter colors denote a decrease in the information value.

Table 7. Potential rules for fluvial channel sand data set.

No	Condition Attribute			Occurrence $C_{j,o}$			Significance $S_{j,o}$		
	O	S	W	K^{low}	K^{med}	K^{high}	K^{low}	K^{med}	K^{high}
1	0	0	0	32	18	0	0.640	0.000	-1.000
2	0	0	1	49	1	0	0.980	-0.980	-1.000
3	0	0	2	50	0	0	1.000	-1.000	-1.000
4	0	1	0	4	46	0	0.000	0.920	-1.000
5	0	1	1	23	26	1	0.000	0.000	-0.980
6	0	1	2	34	16	0	0.680	0.000	-1.000
7	0	2	0	1	49	0	-0.980	0.980	-1.000
8	0	2	1	3	39	8	0.000	0.780	0.000
9	0	2	2	14	30	6	0.000	0.600	0.000
10	1	0	0	2	47	1	0.000	0.940	-0.980
11	1	0	1	23	21	6	0.000	0.000	0.000
12	1	0	2	34	12	4	0.680	0.000	0.000
13	1	1	0	0	50	0	-1.000	1.000	-1.000
14	1	1	1	6	42	2	0.000	0.840	0.000
15	1	1	2	25	23	2	0.500	0.000	0.000
16	1	2	0	5	45	0	0.000	0.900	-1.000
17	1	2	1	6	44	0	0.000	0.880	-1.000
18	1	2	2	15	35	0	0.000	0.700	-1.000
19	2	0	0	0	2	48	-1.000	0.000	0.960
20	2	0	1	0	9	41	-1.000	0.000	0.820
21	2	0	2	0	6	44	-1.000	0.000	0.880
22	2	1	0	0	50	0	-1.000	1.000	-1.000
23	2	1	1	3	46	1	0.000	0.820	-0.980
24	2	1	2	8	34	8	0.000	0.680	0.000
25	2	2	0	29	21	0	0.580	0.000	-1.000
26	2	2	1	35	15	0	0.700	0.000	-1.000
27	2	2	2	42	8	0	0.840	0.000	-1.000

Table 8. Full set of rules sorted according to significance value.

index	<i>O</i>	<i>ca</i> <i>S</i>	<i>W</i>	<i>de</i>	$s_{j,o}$	index	<i>O</i>	<i>ca</i> <i>S</i>	<i>W</i>	<i>de</i>	$s_{j,o}$
19	2	0	0	2	0.262	27	2	2	2	1	-0.52
21	2	0	2	2	0.209	11	1	0	1	2	-0.64
20	2	0	1	2	0.174	14	1	1	1	0	-0.64
3	0	0	2	0	0.112	9	0	2	2	2	-0.64
2	0	0	1	0	0.107	17	1	2	1	0	-0.64
27	2	2	2	0	0.072	21	2	0	2	1	-0.64
22	2	1	0	1	0.068	16	1	2	0	0	-0.7
13	1	1	0	1	0.068	4	0	1	0	0	-0.76
7	0	2	0	1	0.064	12	1	0	2	2	-0.76
10	1	0	0	1	0.058	23	2	1	1	0	-0.82
23	2	1	1	1	0.055	8	0	2	1	0	-0.82
4	0	1	0	1	0.055	15	1	1	2	2	-0.88
16	1	2	0	1	0.052	19	2	0	0	1	-0.88
17	1	2	1	1	0.049	10	1	0	0	0	-0.88
26	2	2	1	0	0.043	14	1	1	1	2	-0.88
14	1	1	1	1	0.043	7	0	2	0	0	-0.94
12	1	0	2	0	0.039	5	0	1	1	2	-0.94
6	0	1	2	0	0.039	10	1	0	0	2	-0.94
8	0	2	1	1	0.035	2	0	0	1	1	-0.94
1	0	0	0	0	0.033	23	2	1	1	2	-0.94
18	1	2	2	1	0.026	25	2	2	0	2	-1.0
25	2	2	0	0	0.024	3	0	0	2	2	-1.0
24	2	1	2	1	0.024	26	2	2	1	2	-1.0
9	0	2	2	1	0.016	21	2	0	2	0	-1.0
15	1	1	2	0	0.014	7	0	2	0	2	-1.0
5	0	1	1	1	0.009	16	1	2	0	2	-1.0
11	1	0	1	0	0.009	13	1	1	0	2	-1.0
5	0	1	1	0	0.009	13	1	1	0	0	-1.0
15	1	1	2	1	0.005	17	1	2	1	2	-1.0
25	2	2	0	1	0.003	19	2	0	0	0	-1.0
11	1	0	1	1	0.003	4	0	1	0	2	-1.0
1	0	0	0	1	0.001	2	0	0	1	2	-1.0
6	0	1	2	1	-0.04	1	0	0	0	2	-1.0
18	1	2	2	0	-0.1	22	2	1	0	2	-1.0
26	2	2	1	1	-0.1	20	2	0	1	0	-1.0
9	0	2	2	0	-0.16	18	1	2	2	2	-1.0
12	1	0	2	1	-0.28	6	0	1	2	2	-1.0
20	2	0	1	1	-0.46	27	2	2	2	2	-1.0
24	2	1	2	2	-0.52	3	0	0	2	1	-1.0
8	0	2	1	2	-0.52	22	2	1	0	0	-1.0
24	2	1	2	0	-0.52						

Table 9. The derived positive rules.

Good O (2)	and	low S (0)	and	No matter W	leads to	high K (2)
Bad O (0)	and	low S (0)	and	large W (2)	leads to	low K (0)
Bad O (0)	and	low S (0)	and	large W (1)	leads to	low K (0)
Good O (2)	and	high S (2)	and	large W (2)	leads to	low K (0)
Good O (2)	and	medium S (1)	and	no matter W	leads to	medium K (1)
Medium O	and	no matter S	and	small W (0)	leads to	medium K (1)
Bad O (0)	and	high S (2)	and	small W (0)	leads to	medium K (1)

Table 10. The derived negative rules.

Good O (2)	and	medium S (1)	and	small W	never leads to	low K (0)
Bad O (0)	and	small S (0)	and	large W (2)	never leads to	low K (0)
Good O (2)	and	high S (2)	and	large W (2)	never leads to	high K (2)
Bad O (0)	and	medium S (1)	and	large W (2)	never leads to	high K (2)
Medium O	and	high S (2)	and	large W (2)	never leads to	high K (2)
Good O (2)	and	small S (0)	and	medium W	never leads to	low K (0)

Table 11. Results of classification.

	1	2	3	accuracy
1	353	90	0	79.68%
2	135	583	17	79.32%
3	12	27	133	77.33%

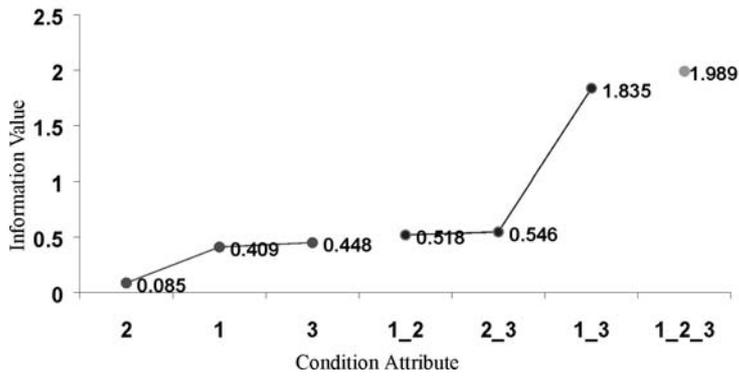


Fig. 6. Information value of all subsets of condition attributes.

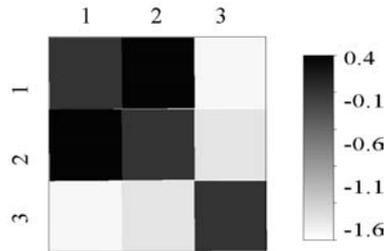


Fig. 7. Information changes for lumped/merged decision categories.

5.2 Facies Assignment from Wireline Well Logs

Theoretically there are $324 (= 3 \times 4 \times 3 \times 3 \times 3)$ configurations of condition attributes; however, only 54 configurations have nonzero coverage. There are $31 (= 2^5 - 1)$ subsets of condition attributes and the complete rule tables are obtained from the entire attribute subset supplemented with “missing” entries filled with rules from the subsets. The top plot of Figure 8 shows the ranked information measure of all 31 subsets of condition attributes. The bottom plot of Figure 8 shows the ranked information measure within groups of subsets with the same number of condition attributes. Even though there is no universal increase in the information measure as the increase of number of condition attributes, the largest information in the subsets with the same number of condition attributes increases as the number of condition attributes in the subsets increases. For this example, the entire set has the largest information measure; therefore the entire set of variables should be used for rule induction.

The complete set of rules and probability table contain 324 lines (configurations). In each line, the positive numbers are probability values for decision outcome with a significant positive *significance* and the negative numbers are the probabilities for decision outcome with a significant negative *significance* added a minus sign. Table 12 lists several lines of the final rule table filtered by the significance threshold and Table 13 lists the same lines in the probability table. The differences between these two tables are the effects of the significance thresholds used for filtering the rules.

The most significant rules have been compared favorably with the empirical rules summarized by geologists who provide the data. Instead of rigid rules summarized by the geologists, for a given set of well logs the rules generated from the proposed method not only provide the most likely facies, but also supplied the probability to be other facies and the information exclusion to be some facies, which would be flexible for the geologist to incorporate their expertise.

The rules were applied to the training data set. Figure 9 presents the rules when applied to the training data in one of the wells. The top plot shows the positive rules which are the probabilities of facies appearing along the depth and the bottom plot shows the negative rules that are the probabilities of facies unlikely to appear.

Figure 10 shows the true cored facies of the training data in Well one (top) and the assigned facies based on well logs (bottom).

For this example, the information changes when *lumping* decision outcome pair ise are shown in Figure 2. The explanation of the distinguishability has already been described previously.

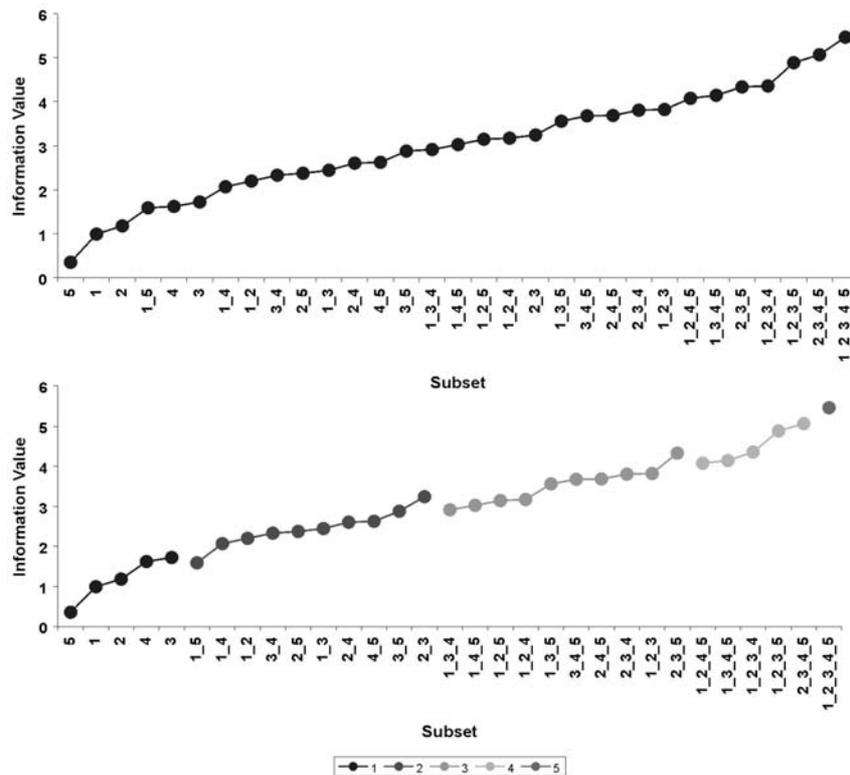


Fig. 8. Information of all subsets of condition attributes, *top*: overall sorted; *bottom*: sorted subset-wisely.

Table 12. Rules in the complete rule table.

<i>No</i>	Code of <i>ca</i>					Filtered Significance of Decision Values									<i>a</i>	<i>b</i>	<i>c</i>	
1	0	0	0	0	0	.00	.00	.93	.00	.00	.00	.00	.00	.00	.00	3	23	15
...	
66	0	2	1	0	2	-1.00	-1.00	-1.00	.00	.00	.00	.00	.22	.22	4	30	36	
67	0	2	1	1	0	1.00	.00	.00	.00	.00	.00	.00	.00	.00	5	31	8	
68	0	2	1	1	1	-1.00	.00	.22	.39	-1.00	.22	-.99	.00	.00	4	30	99	
69	0	2	1	1	2	.00	.00	-1.00	.00	.58	.17	.00	.00	-.99	4	30	98	
70	0	2	1	2	0	1.00	.00	.00	.00	.00	.00	.00	.00	.00	4	27	8	
71	0	2	1	2	1	-1.00	.00	.00	.36	-1.00	.22	.00	.00	.00	3	23		
72	0	2	1	2	2	.00	.00	.00	.00	1.00	.00	.00	.00	.00	4	30	5	
...	
324	2	3	2	2	2	.00	.00	.00	.00	.00	.00	.00	.00	.00	4	28	11	

CA, Condition attribute; *a*, level of subset; *b*, index of subset; *c*, configuration coverage.

Table 13. Conditional probability in the complete probabaility table.

<i>No</i>	Code of <i>ca</i>					Filtered Significance of Decision Values									<i>a</i>	<i>b</i>	<i>c</i>
1	0	0	0	0	0	.00	.00	.93	.07	.00	.00	.00	.00	.00	3	23	15
...
66	0	2	1	0	2	.00	.00	.00	.02	.14	.19	.19	.22	.22	4	30	36
67	0	2	1	1	0	1.00	.00	.00	.00	.00	.00	.00	.00	.00	5	31	8
68	0	2	1	1	1	.00	.11	.22	.39	.00	.22	.01	.02	.02	4	30	99
69	0	2	1	1	2	.03	.06	.00	.06	.58	.17	.06	.02	.01	4	30	98
70	0	2	1	2	0	1.00	.00	.00	.00	.00	.00	.00	.00	.00	4	27	8
71	0	2	1	2	1	.00	.09	.18	.36	.00	.22	.12	.02	.02	3	23	
72	0	2	1	2	2	.00	.00	.00	.00	1.00	.00	.00	.00	.00	4	30	5
...
324	2	3	2	2	2	.09	.27	.00	.27	.36	.00	.00	.00	.00	4	28	11

CA, Condition attribute; *a*, level of subset; *b*, index of subset; *c*, configuration coverage

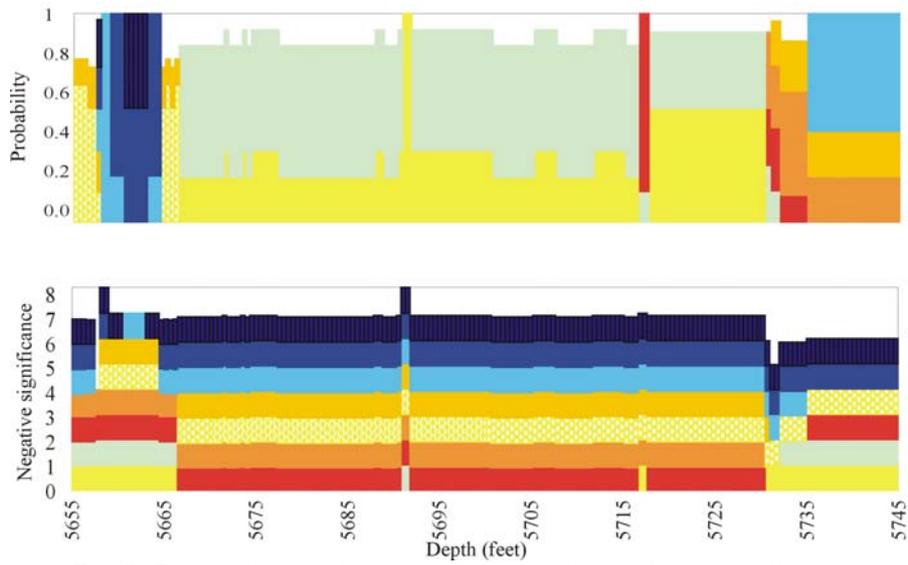


Fig. 9. Positive (top) and negative (bottom) rules applying to well one.

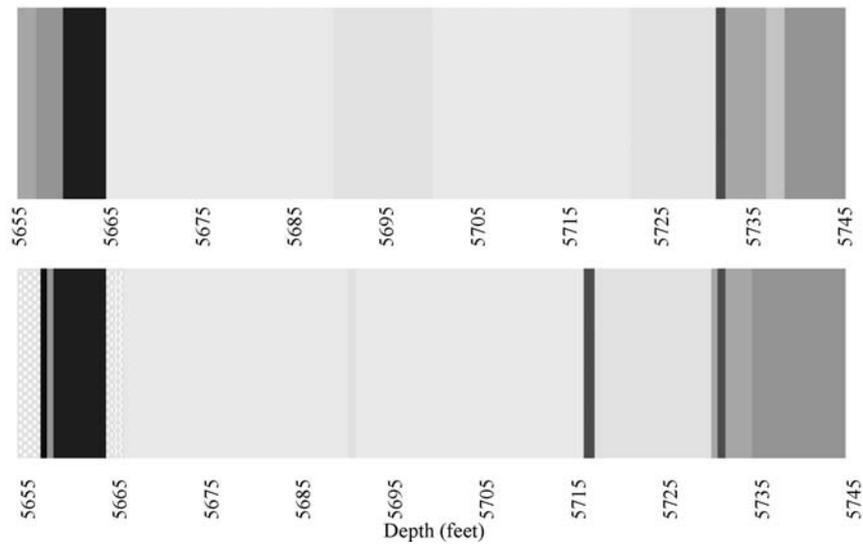


Fig. 10. True facies (top) and assigned facies based on the rules (bottom) of well one.

5.3 Conclusion

The proposed technique has been successfully applied in the two example data sets. Some significant rules have been retrieved from the fluvial system example which are actually embedded in the data generation. It is understandable to have high effective permeability if the channels and channel complex have a good direction, low sinuosity. Bad direction and low sinuosity or good direction with high sinuosity lead to low effective permeability. The rules for the facies assignment from well logs are the facies to be assigned for a given set of well log values. The rule compared favorably with the ones derived by the geologists working extensively in the field where data was collected. The rules provide not only the most likely facies for a given set of well log values, but also provide the probability to be other facies. The results from the distinguish analysis actually supplies hints to the geologist to re-consider the distinction of some facies.

5.4 Discussion

The proposed algorithm works well for the examples. The most important positive and negative rules are retrieved successfully. The proposed *significance* definition combines measures of *accuracy* and *coverage* and serves as a quality measure of rules. Also, the significance identifies positive and negative rules, similar to the positive region and negative region in rough sets.

The proposed rule induction technique is suited to geological data where most attributes are significant. The proposed significance measure can be used in combination with other rule induction techniques and serves as a ranking measure to identify the most important rules. In general, however, the algorithm will need to be extended to include attribute reduction.

Careful examination of a data table may lead an experienced person to infer similar, if not the same rules; however, there are many advantages to automatic rule induction. The procedure works for very large datatables with many attributes, it is repeatable, and avoids personal biases. It could also lead to nonintuitive, but meaningful data relations. The effort in making predictions can be greatly reduced.

Glossary

Accuracy: a measure of the closeness of a conditional probability to 1 or 0.

Condition attribute value: a measure of the value contributed by every possible combination of conditioning attributes to a decision outcome.

Conditional probability: for each decision outcome and condition configuration pair, the fraction of occurrences in the data table out of configuration coverage.

Conditional probability table: a complete table of conditional probability for all decision outcome and all condition configuration pairs.

Configuration: a unique combination of the values taken by the condition attributes.

Configuration coverage: the number of observations associated with each configuration.

Data table: table of observed decisions for each instance of the combination of conditioning attributes. The set of conditioning attributes are analyzed through basic statistics and clustering so that class assignments can be made.

Decision coverage: number of observations associated with each decision outcome.

Distinguishability: a measure of distinctiveness between decision outcomes, based on the expected prediction probabilities.

Relative configuration coverage: a measure of closeness to sufficiency, defined as a proportionality of occurrence frequency of a configuration and decision outcome pair related to the configuration coverage.

Relative decision coverage: a measure of the proportion of a configuration and decision outcome pair out of the number for that outcome.

Rule table: a table of the positive and negative rules with associated conditional probabilities, which are filtered by significance.

Significance: a measure of how an outcome is both sufficient and accurate; in practice, a combination of the relative coverage, accuracy, and decision coverage that leads to a rule.

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