

Machine Learning Modelling in R : : CHEAT SHEET

Supervised & Unsupervised Learning

Meta-Algorithm, Time Series & Model Validation

| | ALGORITHM | DESCRIPTION | R PACKAGE::FUNCTION | SAMPLE CODE | | ALGORITHM | DESCRIPTION | R PACKAGE::FUNCTION | SAMPLE CODE |
|-----------------------|--|--|--|--|------------------|--|---|--|--|
| SUPERVISED LEARNING | NBC Naïve Bayes classifier | A classification technique based on Bayes' Theorem with an assumption of independence among predictors. In simple terms, a Naïve Bayes classifier assumes that the presence of a particular feature in a class is unrelated to the presence of any other feature | e1071::naiveBayes | naiveBayes(class ~, data = x) | META ALGORITHM | REGU Regularisation L1 (Lasso) L2 (Ridge) | Regularization adds a penalty on the different parameters of a model to reduce the freedom of the model. Hence, the model will be less likely to fit the noise of the training data and will improve the generalization abilities of the model | glmnet::glmnet | L1 : glmnet(myMatrixA, myMatrixB, family = "gaussian", alpha = 1) L2 : glmnet(myMatrixA, myMatrixB, family = "gaussian", alpha = 0) |
| | KNN k-Nearest Neighbours | A non-parametric method used for classification and regression. In both cases, the input consists of the k closest training examples in the feature space. The output depends on whether k-NN is used for classification or regression | class::knn | knn(train, test, cl, k = 1, l = 0, prob = FALSE, use.all = TRUE) | | BOO Boosting | A process of iteratively refining, e.g. by reweighting, of estimated regression and classification functions (though it has primarily been applied to the latter), in order to improve predictive ability. | Parametric model rpart::rpart | glmboost(Yen ~, data = curr1(trndids,)) |
| | REG Linear Regression | Model the linear relationship between a scalar dependent variable Y and one or more explanatory variables (or independent variables) denoted X | stats::lm | lm(dist ~ speed, data=cars) | | BAG Bagging | Bagging is a way to increase the power of a predictive statistical model by taking multiple random samples (with replacement) of the training data set, and using each of them to construct a separate model and separate predictions for the original test set | All models: foreach Tree models: ipred::bagging | foreach : d <- data.frame(x=1:10, y=rnorm(10)) s <- foreach(d=iter(d, by='row', combine='bind')) %doParallel %dopar { identical(s, d) } ipred: bagging(formula, data, subset, na.action=na.rpart, ...) |
| | LOG Logistic Regression | Used to predict a binary outcome (1 / 0, Yes / No, True / False) given a set of independent variables. | stats::glm | glm(Y ~, family = binomial (link = 'logit'), data = X) | | PRU Pruning | Pruning is a technique that reduces the size of decision tree by removing sections of the tree that provide little power to classify instances. Pruning reduces the complexity of the final classifier and hence improves predictive accuracy by reducing overfitting | rpart::prune | prune(x, cp = 0.1) |
| | TM Tree-Based Models | The idea is to consecutively divide (branch) the training dataset based on the input features until an assignment criterion with respect to the target variable into a "data bucket" (leaf) is reached | rpart::rpart | rpart(Kyphosis ~ Age + Number + Start, data = kyphosis) | | RFO Random Forest | An ensemble learning method for classification, regression and other tasks, that operate by constructing a multitude of decision trees at training time and outputting the class that is the mode of the classes (classification) or mean prediction (regression) | randomForest::randomForest | randomForest(X ~, data = Y, subset = mySub) |
| UNSUPERVISED LEARNING | ANN Artificial Neural Network | Neural networks are built from units called perceptrons. Perceptrons have one or more inputs, an activation function and an output. An ANN model is built up by combining perceptrons in structured layers. | neuralnet::neuralnet | neuralnet(f, data=train_hidden=c(5,3), linear.output=T) | MODEL VALIDATION | STS Lead-lag analysis, Auto-correlation, Spectral analysis, Time series clustering, Seasonality, Trend,... | Random sampling of observations for training and testing a model can be an issue when faced with a times dimension. Random sampling may either destroy serial correlation properties in the data which we would like to exploit | stats forecast spectral TTR | Auto-correlation : acf(x, lag.max = NULL, type = c("correlation", "covariance", "partial")) Spectral Analysis : spec.pgram(myTs, spans = NULL) Seasonal Decomposition of Time Series : stl(x, window = 7, l.window = 50, l.jump = 1) |
| | SVM Support Vector Machine | A data classification method that separates data using hyperplanes | e1071::svm | svm(formula, data = NULL, ..., subset, na.action = na.omit, scale = TRUE) | | PM Performance metrics | Depends on the problem: • Regression : squared errors, outliers, error rate... • Classification : Accuracy, precision, recall, F-score... | Regression : stats::outlierTest, stats::qqPlot... Classification-ROC-R: Tree : caret::confusionMatrix | Regression : fit <- lm(Y~X, data=myData) outlierTest(fit) qqPlot(fit, main="QQ Plot") |
| | PCA Principal Component Analysis | A procedure that uses an orthogonal transformation to convert a set of observations of possibly correlated variables into a set of values of linearly uncorrelated variables called principal components. | stats::princomp stats::princomp FactoMineR::PCA ade4::dudi.pca amap::acp | stats : princomp(formula, data = NULL, subset, na.action, ...) FactoMineR : PCA(decathlon, quanti.sup = 11:12, quali.sup = 13) ade4 : dudi.pca(deugStab, center = deugScnt, scale = FALSE, scan = FALSE) amap : acpl(busch) | | BVT Bias-Variance Tradeoff | Simple models with few parameters are easier to compute but may lead to poorer fits (high bias). Complex models may provide more accurate fits but may over-fit the data (high variance) | Tailored to the analysis | Tailored to the analysis |
| | KMC k-Mean Clustering | Aims at partitioning n observations into k clusters in which each observation belongs to the cluster with the nearest mean | stats::kmeans | kmeans(x, centers, iter.max = 10, nstart = 1, algorithm = c("Hartigan-Wong", "Lloyd", "Forgy", "MacQueen"), trace=FALSE) | | CV Cross validation | Cross validation compares the test performances of different model realisations with different sets or values of parameters | caret::createDataPartition caret::createFolds | createDataPartition(classes, p = 0.8, list = FALSE) |
| | HCL Hierarchical Clustering | An approach which builds a hierarchy from the bottom-up, and doesn't require the number of clusters to be specified beforehand. | stats::hclust | hclust(d, method = "complete", members = NULL) | | LC Learning Curves | Learning curves plot a model's training and test errors, or the chosen performance metric, depending on the training set size | caret::learning_curve_dat | learning_curve_dat(dat, outcome = NULL, proportion = (1:10)/10, test_prop = 0, verbose = TRUE, ...) |

Standard Modelling Workflow

Time Series View

