

# Deep Learning & AutoML in h2o



## Set up

```
# Load libraries
library(h2o)
library(tidyverse)
library(wesanderson)
library(knitr)

# Disable progress bar in document
h2o.no_progress()

# Start h2o cluster
h2o.init(nthreads = -1,
        max_mem_size = '4G')
```

```
## Connection successful!
##
## R is connected to the H2O cluster:
##   H2O cluster uptime:      3 hours 1 seconds
##   H2O cluster timezone:    Europe/Warsaw
##   H2O data parsing timezone: UTC
##   H2O cluster version:     3.40.0.1
##   H2O cluster version age:  1 month and 6 days
##   H2O cluster name:        H2O_started_from_R_User_acc441
##   H2O cluster total nodes: 1
##   H2O cluster total memory: 3.68 GB
##   H2O cluster total cores: 16
##   H2O cluster allowed cores: 16
##   H2O cluster healthy:     TRUE
##   H2O Connection ip:       localhost
##   H2O Connection port:     54321
##   H2O Connection proxy:    NA
##   H2O Internal Security:   FALSE
##   R Version:                R version 4.2.2 (2022-10-31 ucrt)
```

```

# Load file
mushrooms <- read.csv('https://tinyurl.com/hmkhs9au')

# Transform data set for further analysis
mushrooms <- mushrooms %>%
  # Remove not needed characters
  mutate(across(1:23, ~ substr(.x, 3,3))) %>%
  # Change columns from strings to factors
  mutate(across(everything(), as.factor))

# Load data to h2o cluster
mushrooms_hex <- as.h2o(mushrooms, destination_frame = 'mushrooms_hex')

# Split data set to train (75%) & test (25%)
mushrooms_split <- h2o.splitFrame(data = mushrooms_hex, ratios = 0.75)

mushrooms_train <- mushrooms_split[[1]]
mushrooms_test <- mushrooms_split[[2]]

```

## Data

Data set contains 8124 observations of 23 species from Agaricus and Lepiota families. Beside class (e - edible/p - poisonous or unknown edibility) there are 22 physical attributes.

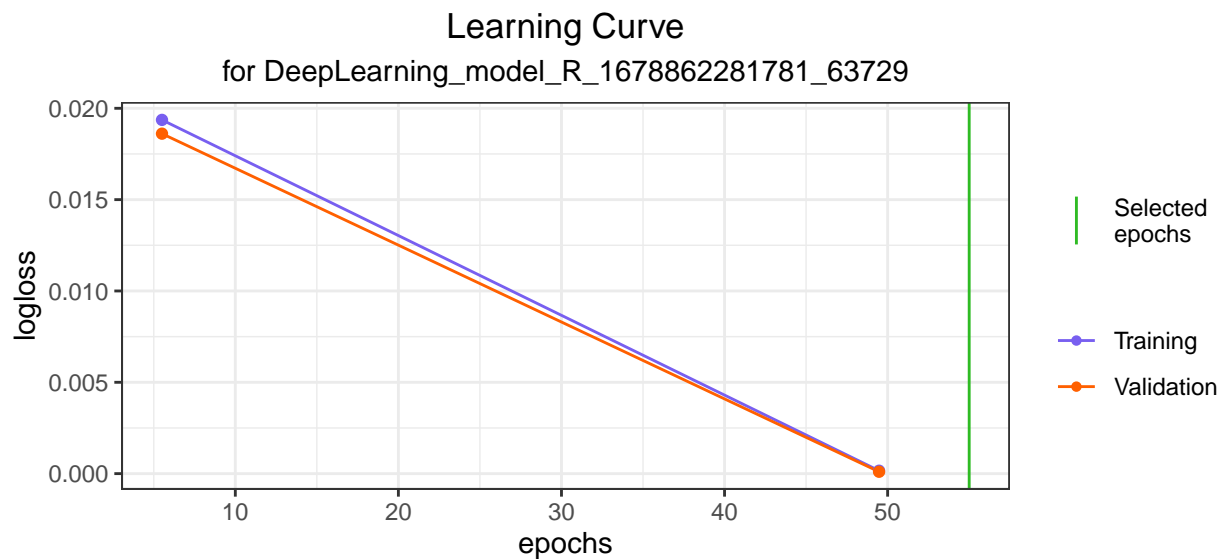
## h2o's Deep Learning

```

# Build and train Deep Learning model
mushrooms_dl <- h2o.deeplearning(
  y = 23,
  x = 1:22,
  training_frame = mushrooms_train,
  validation_frame = mushrooms_test,
  distribution = 'multinomial',
  activation = 'RectifierWithDropout',
  hidden = c(100, 200, 100),
  input_dropout_ratio = 0.2,
  l1 = 1e-5,
  epochs = 55,
  variable_importances = TRUE)

```

```
# Learning curve plot
h2o.learning_curve_plot(mushrooms_dl)
```



```
# Confusion matrix
h2o.confusionMatrix(mushrooms_dl, mushrooms_test, valid = FALSE, xval = FALSE)
```

```
## Confusion Matrix (vertical: actual; across: predicted) for max f1 @ threshold = 0.86974572554556:
##      e    p  Error   Rate
## e   1016    0 0.000000 =0/1016
## p      0 1004 0.000000 =0/1004
## Totals 1016 1004 0.000000 =0/2020
```

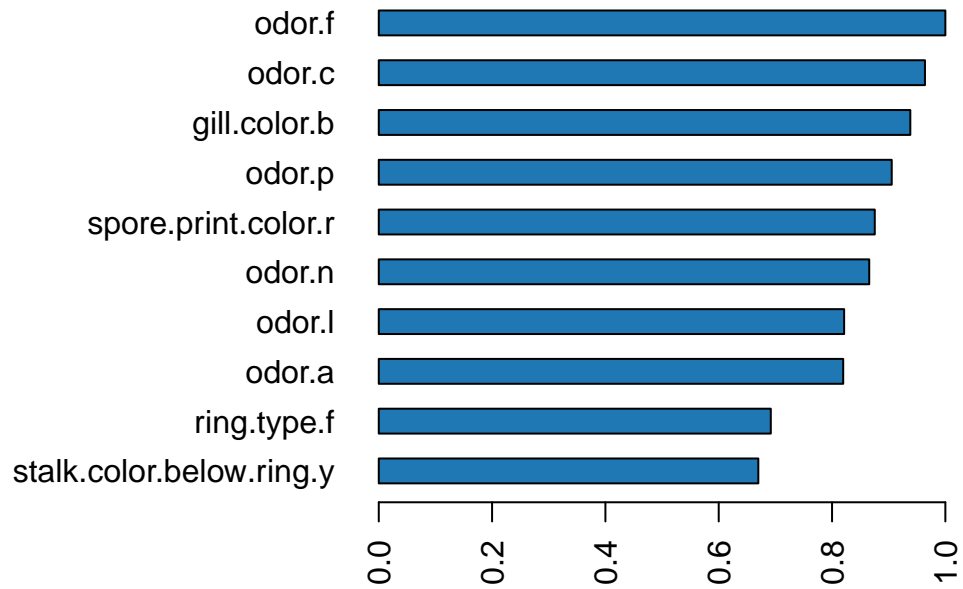
```
# If needed, predictions can be done on test data set
# (which was used for validation in model training)
mushroom_pred <- h2o.predict(object = mushrooms_dl,
                             newdata = mushrooms_test)

# Confusion matrix as simple table: test vs predictions
table(as.data.frame(mushrooms_test[,23])[,1],
      as.data.frame(mushroom_pred[,1])[,1])
```

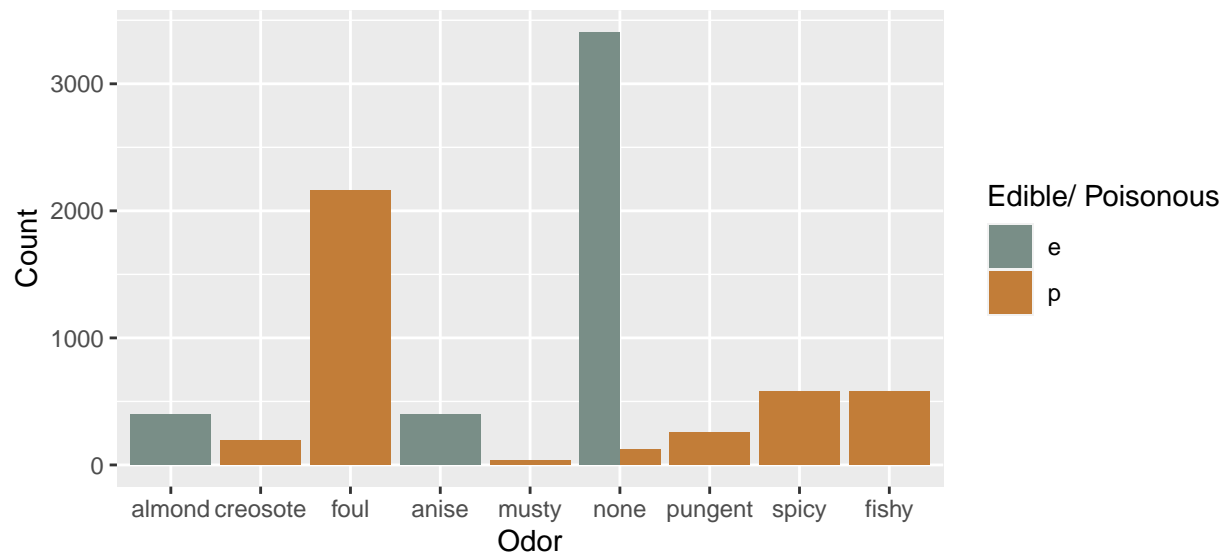
```
##
##      e    p
## e 1016    0
## p    0 1004
```

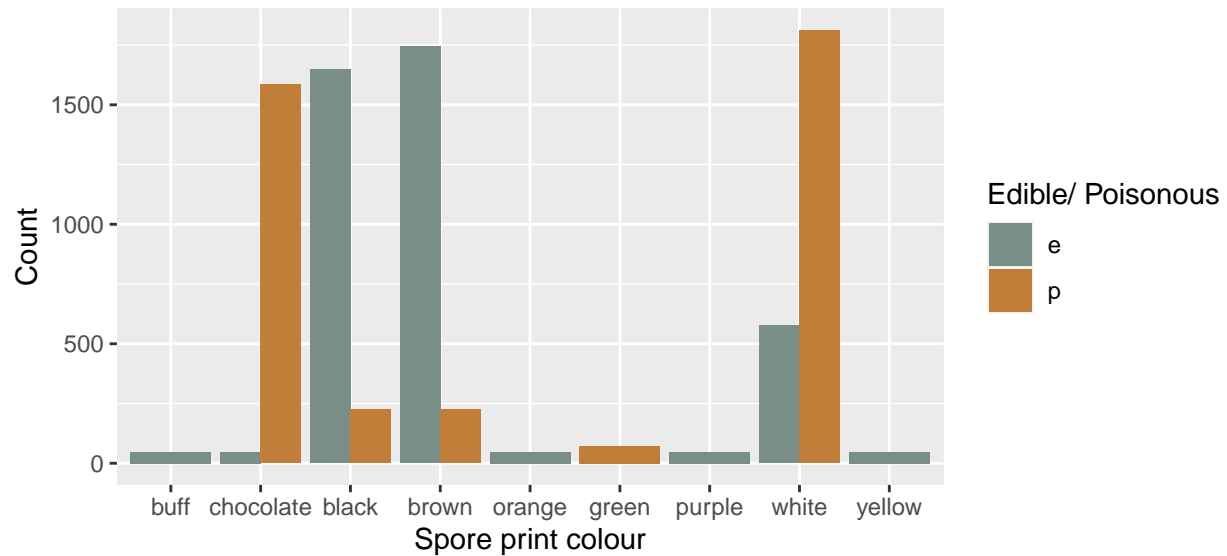
```
# Importance of parameters
h2o.varimp_plot(mushrooms_dl)
```

### Variable Importance: Deep Learning



Most important parameters are odor and spore print color:





## AutoML

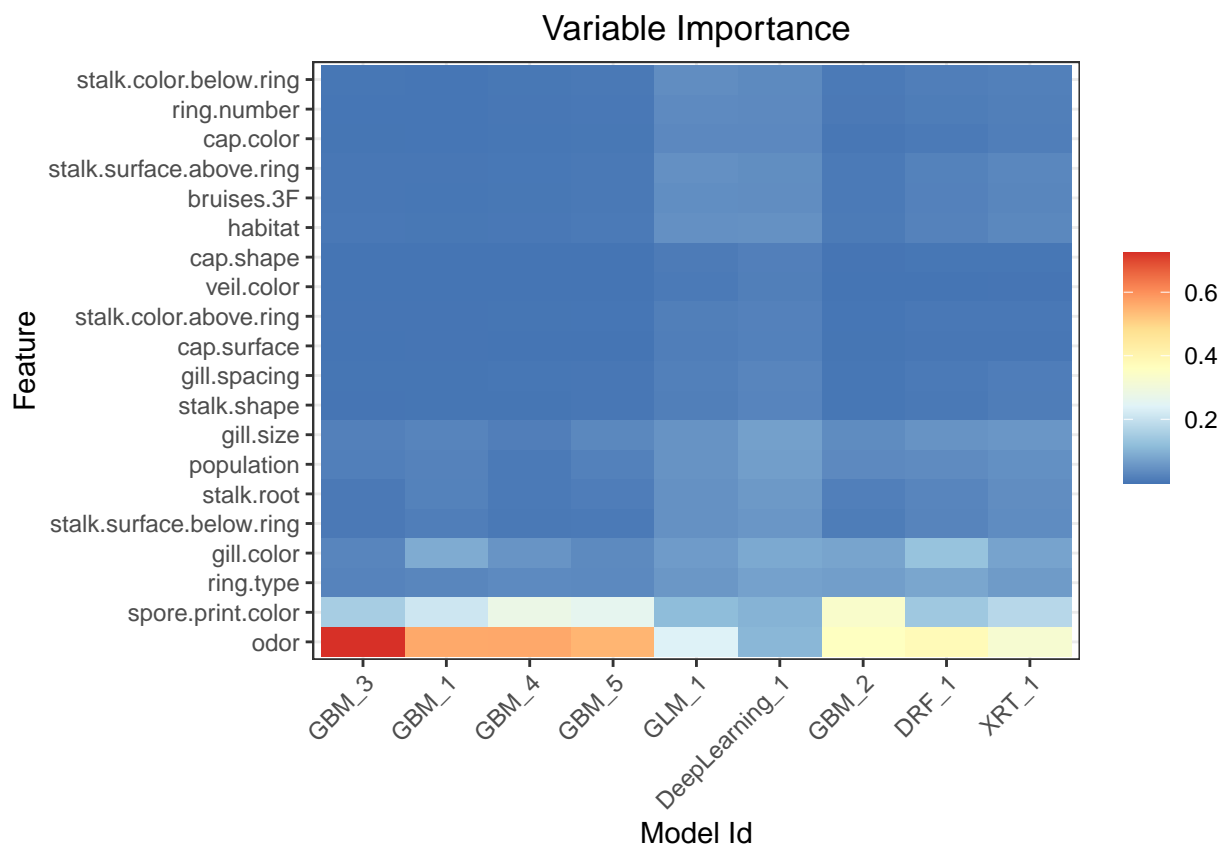
```
# Run AutoML: max 100 models, max 60 seconds
mushrooms_auoml <- h2o.automl(
  y = 23,
  x = 1:22,
  training_frame = mushrooms_train,
  max_runtime_secs = 60,
  max_models = 100)

##
## 10:38:26.963: AutoML: XGBoost is not available; skipping it.
## 10:38:26.968: _train param, Dropping bad and constant columns: [veil.type]
## 10:38:27.481: _train param, Dropping bad and constant columns: [veil.type]
## 10:38:40.666: _train param, Dropping bad and constant columns: [veil.type]
## 10:38:41.55: _train param, Dropping bad and constant columns: [veil.type]
## 10:38:48.666: _train param, Dropping bad and constant columns: [veil.type]
## 10:38:56.890: _train param, Dropping bad and constant columns: [veil.type]
## 10:39:05.553: _train param, Dropping bad and constant columns: [veil.type]
## 10:39:06.108: _train param, Dropping bad and constant columns: [veil.type]
## 10:39:12.414: _train param, Dropping bad and constant columns: [veil.type]

# Leader board
df <- h2o.get_leaderboard(object = mushrooms_auoml, extra_columns = "ALL")
df <- as.data.frame(df)
df$model_id <- substr(df$model_id,1,5)
df <- df %>% mutate(across(where(is.numeric), round, 3))
df <- df[,-10]
kable(df,
  col.names = c('Model', 'auc', 'Log Loss', 'aucPR', 'Mean per class err.', 'RMSE',
    'MSE', 'Train. time [ms]', 'Predict time / row [ms]'))
```

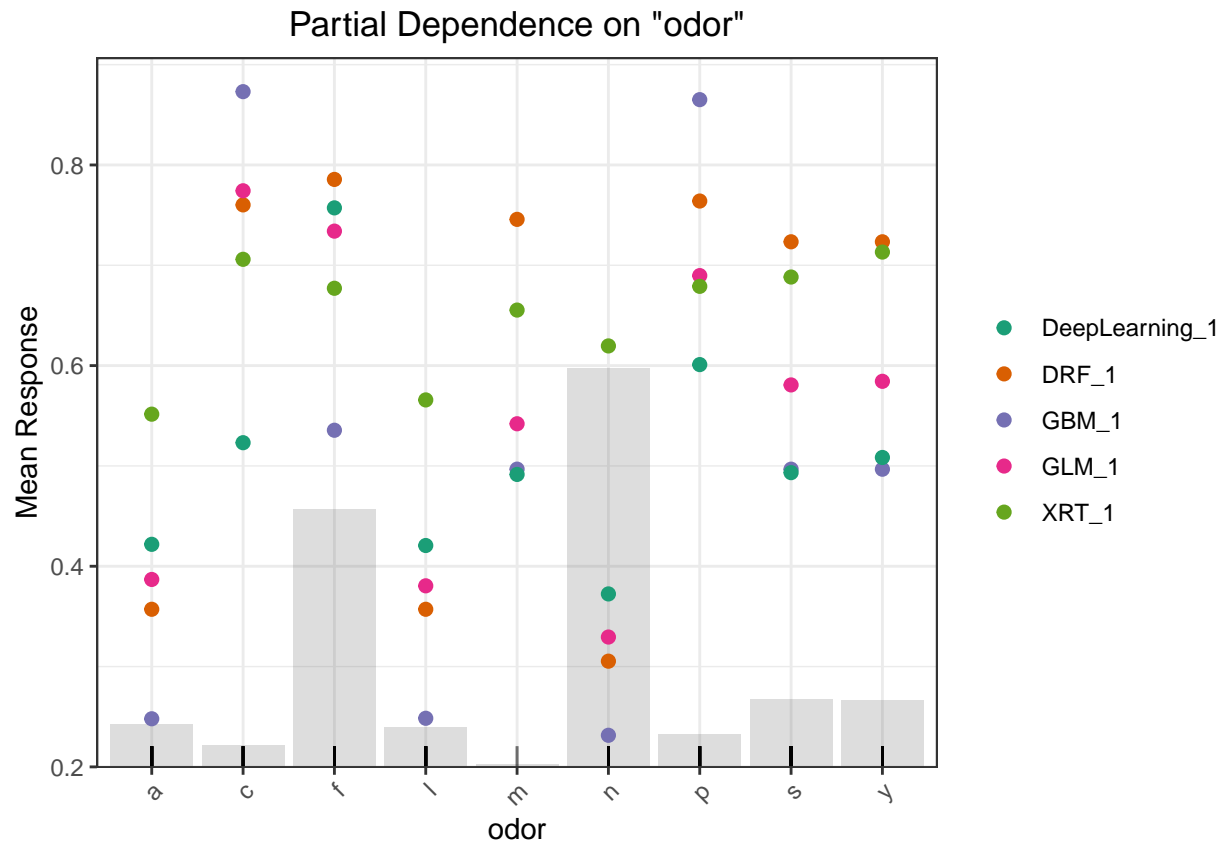
Model	auc	Log Loss	aucPR	Mean per class err.	RMSE	MSE	Train. time [ms]	Predict time / row [ms]
DRF_1	1	0.005	1	0.000	0.013	0.000	91	0.004
GBM_1	1	0.000	1	0.000	0.000	0.000	3587	0.116
GBM_2	1	0.000	1	0.000	0.000	0.000	2069	0.065
GBM_4	1	0.000	1	0.000	0.000	0.000	2318	0.074
GBM_3	1	0.000	1	0.000	0.000	0.000	2144	0.069
GBM_5	1	0.000	1	0.000	0.008	0.000	1623	0.055
GLM_1	1	0.001	1	0.000	0.007	0.000	110	0.002
DeepL	1	0.001	1	0.000	0.018	0.000	224	0.005
XRT_1	1	0.313	1	0.001	0.301	0.091	115	0.005

```
# Variables importance heatmap for different AutoML models
h2o.varimp_heatmap(mushrooms_auml)
```



```
# Effect of odor variable for each model
```

```
h2o.pd_multi_plot(mushrooms_auoml, mushrooms_test, "odor")
```



## References

1. **Data set:** <https://www.kaggle.com/datasets/ulrikthgepedersen/mushroom-attributes>
2. **Agaricus family graphic:** <https://en.wikipedia.org/wiki/Agaricus>
3. **Mushrooms graphics:** <<https://biolwww.usask.ca/fungi/glossary.html>>
4. **h2o Deep Learning:** <https://docs.h2o.ai/h2o/latest-stable/h2o-docs/data-science/deep-learning.html>
5. **h2o AutoML:** <https://docs.h2o.ai/h2o/latest-stable/h2o-docs/automl.html>
6. **h2o models explainability:** <https://docs.h2o.ai/h2o/latest-stable/h2o-docs/explain.html#explanation-plotting-functions>