

# pdp: An R Package for Creating Partial Dependence Plots

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**Abstract** Complex nonparametric models—like neural networks, random forests, and support vector machines—are more common than ever in predictive analytics, especially when dealing with large observational databases that don't adhere to the strict assumptions imposed by traditional statistical techniques (e.g., multiple linear regression). Unfortunately, it can be challenging for the uninitiated to understand the results of such models and explain them to management. Partial dependence plots offer a simple solution. Partial dependence plots are low-dimensional graphical renderings of the prediction function  $\hat{f}(x)$  so that the relationship between the outcome and predictors of interest can be easily understood. These plots are especially useful in explaining the output from black box models. In this paper, we introduce **pdp**, a general R package for constructing partial dependence plots.

## Introduction

Predictor importance is an important task in any supervised learning problem. However, ranking variables is only part of the story and once a subset of "important" features is identified it is often necessary to assess the relationship between them (or subset thereof) and the response. This can be done in many ways, but in machine learning it is often accomplished by constructing *partial dependence plots* (PDPs); see Friedman (2000) for details. PDPs help visualize the relationship between a subset of the predictors (typically 1-3) and the response while accounting for the average effect of the other independent variables in the model. They are particularly effective with black box models like random forests and support vector machines.

Let  $x = \{x_1, x_2, \dots, x_p\}$  represent the predictors in a model whose prediction function is  $\hat{f}(x)$ . If we partition  $x$  into an interest set,  $z_s$ , and its complement,  $z_c = x \setminus z_s$ , then the "partial dependence" of the response on  $z_s$  is defined as

$$f_s(z_s) = E_{z_c} [\hat{f}(z_s, z_c)] = \int \hat{f}(z_s, z_c) p_c(z_c) dz_c, \quad (1)$$

where  $p_c(z_c)$  is the marginal probability density of  $z_c$ :  $p_c(z_c) = \int p(x) dz_s$ . Equation (1) can be estimated from a set of training data by

$$\bar{f}_s(z_s) = \frac{1}{n} \sum_{i=1}^n \hat{f}(z_s, z_{i,c}), \quad (2)$$

where  $z_{i,c}$  ( $i = 1, 2, \dots, n$ ) are the values of  $z_c$  that occur in the training sample; that is, we average out the effects of all the other predictors in the model.

Constructing a PDP (2) in practice is rather straightforward. To simplify, let  $z_s = x_1$  be the predictor variable of interest with unique values  $\{x_{11}, x_{12}, \dots, x_{1k}\}$ . The partial dependence of the response on  $x_1$  can be constructed as follows:

1. For  $i \in \{1, 2, \dots, k\}$ :
  - (a) Create a copy of the training data and replace the original values of  $x_1$  with the constant  $x_{1i}$ .
  - (b) Compute the vector of predicted values from the modified copy of the training data.
  - (c) Compute the average prediction to obtain  $\bar{f}_1(x_{1i})$ .
2. Plot the pairs  $[x_{1i}, \bar{f}_1(x_{1i})]$  for  $i = 1, 2, \dots, k$ .

This can be quite computationally intensive since the algorithm involves  $k$  passes over the training records. Fortunately, this can be parallelized quite easily (more on this in a later section). This algorithm can also be easily extended to larger subsets of two or more features as well.

Limited implementations of Friedman's PDPs are available in packages **randomForest** (Liaw and Wiener, 2002) and **gbm**, among others; these are limited in the sense that they only apply to the models fit using the respective package. For example, the `partialPlot` function in **randomForest** only applies to objects of class "randomForest". While the **randomForest** implementation will only allow for a single predictor, the **gbm** implementation can deal with any subset of the predictor space. Partial dependence functions are not restricted to tree-based models; they can be applied to any supervised

learning algorithm (e.g., multiple linear regression and neural networks). However, to our knowledge, there is no general package for constructing PDPs in R (e.g., PDPs for a *conditional random forest* as implemented by the `cforest` function in the `party` and `partykit` packages; see [Torsten Hothorn and Zeileis \(2015\)](#) and [Hothorn and Zeileis \(2016\)](#), respectively). The `pdp` ([Greenwell, 2016](#)) package tries to close this gap by offering a general framework for constructing PDPs that can be applied to several types of fitted models.

The `plotmo` package ([Milborrow, 2015](#)) is one alternative to `pdp`. According to the author, `plotmo` constructs "a poor man's partial dependence plot." In particular, it plots a model's response when varying one or two predictors while holding the other predictors in the model constant (continuous features are fixed at their median value, while factors are held at their first level). These plots allow for up to two variables at a time and are less accurate than PDPs, but are faster to construct. For additive models (i.e., models with no interactions), these plots are identical in shape to PDPs.

PDPs can be misleading in the presence of substantial interactions ([Goldstein et al., 2015](#)). To overcome this issue [Goldstein, Kapelner, Bleich, and Pitkin](#) developed the concept of *individual conditional expectation* (ICE) plots—available in the `ICEbox` package. `ICEbox` only allows for one variable at a time (i.e., no multivariate displays), though color can be used effectively to display information about an additional predictor. The ability to construct derivative plots is also available.

Many other techniques exist for visualizing relationships between the predictors and the response based on a fitted model. For example, the `car` package ([Fox and Weisberg, 2011](#)) contains many functions for constructing partial-residual and marginal-model plots. The `effects` package ([Fox, 2003](#)) is also of interest. However, these methods are orientated towards simpler parametric models (e.g., linear and generalized linear models), whereas `plotmo`, `ICEbox`, and `pdp` are mainly for nonparametric and black box models (though they can be used for simple parametric models as well).

## Constructing PDPs in R

As described in the introduction, the `pdp` package is useful for constructing PDPs for many types of fitted models in R. These are especially useful for visualizing the relationships discovered by complex machine learning algorithms such as a random forest. The latest stable release is available from [CRAN](#):

```
install.packages("pdp")
```

The development version is located on GitHub: <https://github.com/bgreenwell/pdp>. Bug reports and suggestions are appreciated and should be submitted to <https://github.com/bgreenwell/pdp/issues>.

Currently, only two functions are exported by `pdp`:

- `partial`
- `plotPartial`

The `partial` function evaluates the partial dependence (2) from a fitted model over a grid of predictor values; the predictors are specified using the `pred.var` argument. If `plot = FALSE` (the default), `partial` returns an object of class "partial" which inherits from the class "data.frame"; put another way, by default, `partial` returns a data frame with an additional class that is recognized by the `plotPartial` function. The columns of the data frame are labeled in the same order as the features supplied to `pred.var`, and the last column is always labeled `y` and contains the values of the partial dependence function  $\hat{f}_s(z_s)$ . If `plot = TRUE`, then `partial` makes an internal call to `plotPartial` and returns the PDP in the form of a `lattice` plot ([Sarkar, 2008](#)). The `plotPartial` function can be used for displaying more advanced PDPs; it operates on objects of class "partial" and has many useful plotting options. Of course, since the default output produced by `partial` is still a data frame, the user can use any plotting package he/she desires to visualize the results—`ggplot2` ([Wickham, 2009](#)), for instance.

Currently supported models are described in Table 1. However, the default method `partial.default` is flexible enough to handle many of the model types not listed in in this table; for example, neural networks from the `nnet` package ([Venables and Ripley, 2002](#)) and projection pursuit regression ([Friedman and Stuetzle, 1981](#)) using the `ppr` function in the `stats` package. Although, for these cases, the user will need to specify the additional options `type` (i.e., "regression" or "classification") and `train` (the original training data) in the call to `partial`—the exception in Table 1 is using `partial` with the "xgb.Booster" class, which requires the user to specify `train`.

For illustration, we will use the (corrected) Boston housing data which are available from the `mlbench` package ([Dimitriadou, 2010](#)). These data contain the median value of owner-occupied homes in 506 U.S. census tracts in the Boston area, along with 13 independent variables such as the per capita crime rate by town. We begin by loading the data and omitting unimportant columns.

Type of model	R package	Object class
Decision tree	<b>C50</b> (Kuhn et al., 2015) <b>party</b> <b>partykit</b> <b>rpart</b> (Therneau et al., 2015)	"C5.0" "BinaryTree" "constparty"/"party" "rpart"
Bagged decision trees	<b>adabag</b> (Alfaro et al., 2013) <b>ipred</b> (Peters and Hothorn, 2015)	"bagging" "classbagg", "regbagg"
Boosted decision trees	<b>adabag</b> (Alfaro et al., 2013) <b>gbm</b> <b>xgboost</b>	"boosting" "gbm" "xgb.Booster"
Cubist	<b>Cubist</b> (Kuhn et al., 2014)	"cubist"
Random forest	<b>randomForest</b>	"randomForest"
Conditional random forest	<b>party</b>  <b>partykit</b>	"RandomForest"  "cforest"
Linear model	<b>stats</b>	"lm"
Generalized linear model	<b>stats</b>	"glm", "lme4"
Nonlinear least squares	<b>stats</b>	"nls"
Multivariate adaptive regression splines (MARS)	<b>earth</b> (Milborrow, 2016)	"earth"
Support vector machine	<b>e1071</b> (Meyer et al., 2015) <b>kernlab</b> (Karatzoglou et al., 2004)	"svm" "ksvm"

**Table 1:** Models specifically supported by the **pdp** package. **Note:** package **xgboost** is only partially supported; see Section 2.2.6 for details.

```
data(BostonHousing2, package = "mlbench") # mlbench must be installed!
boston <- BostonHousing2[, -c(1, 2, 5)]
```

Next, we fit a random forest to the entire data set with default tuning parameters and 500 trees:

```
library(randomForest)
set.seed(101) # for reproducibility
boston.rf <- randomForest(cmedv ~ ., data = boston, importance = TRUE)
varImpPlot(boston.rf) # variable importance plot
```

The model fit is reasonable, with an *out-of-bag* (pseudo)  $R^2$  of 0.89. The variable importance scores are displayed in Figure 1. Both plots indicate that the percentage of lower status of the population (*lstat*) and the average number of rooms per dwelling (*rm*) are highly associated with the median value of owner-occupied homes (*cmedv*). The question then arises, "What is the nature of these associations?" To help answer this, we can look at the partial dependence of *cmedv* on *lstat* and *rm*, both individually and together.

### Single predictor PDPs

As previously mentioned, the **randomForest** package has its own **partialPlot** function for visualizing the partial dependence of the response on a single predictor. For example, the following snippet of code plots the partial dependence of *cmedv* on *lstat*:

```
partialPlot(boston.rf, pred.data = boston, x.var = "lstat")
```

The same plot can be achieved using the **partial** function and setting **plot = TRUE** (see the left side of Figure 2):

```
library(partial)
partial(boston.rf, pred.var = "lstat", plot = TRUE)
```

The only difference is that **pdp** uses the **lattice** graphics package to produce all of its displays.

For a more customizable plot, we can set **plot = FALSE** in the call to **partial** and then use the **plotPartial** function on the resulting data frame. This is illustrated in the example below which increases the line width, adds a LOESS smoother, and customizes the *y*-axis label. The result is displayed in the right side of Figure 2. **Note:** the *pipe* operator **%>%** provided by the **magrittr** package (Bache and Wickham, 2014) is exported whenever **pdp** is loaded for writing more readable code.



**Figure 1:** Dotchart of variable importance scores for the Boston housing data based on a random forest with 500 trees.

```
boston.rf %>% # the %>% operator is read as "and then"
  partial(pred.var = "lstat") %>%
  plotPartial(smooth = TRUE, lwd = 2, ylab = expression(f(lstat)))
```

### Multi-predictor PDPs

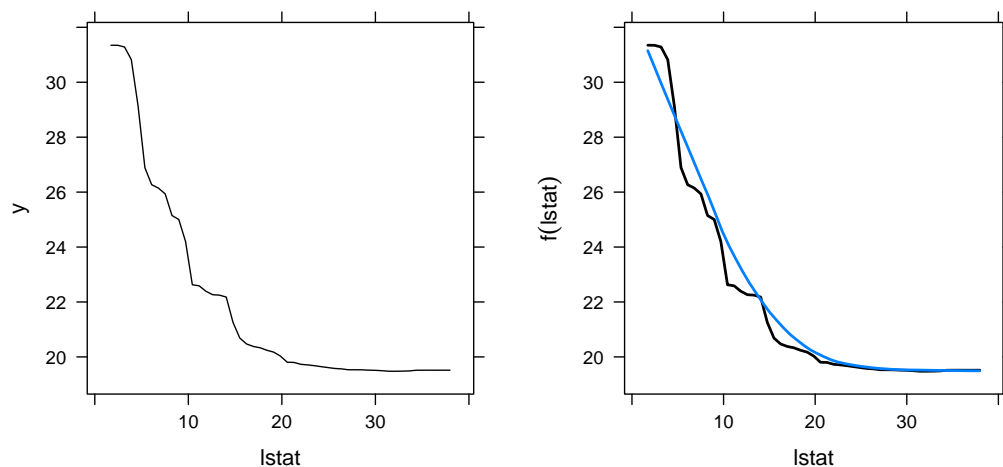
The benefit of using `partial` is threefold: (1) it is a generic function that can be used for various types of fitted models (not just random forests), (2) it will allow for any number of predictors to be used (e.g., multivariate displays), and (3) it can utilize any of the parallel backends supported by the `foreach` package (Analytics and Weston, 2015c); we discuss parallel execution in a later section. For example, the following code chunk uses a multivariate adaptive regression splines (MARS) model (Friedman, 1991) to assess the joint effect of `lstat` and `rm` on `cmedv`. The results, which make use of various `plotPartial` options, are displayed in Figure 3.

```
library(earth) # for fitting MARS models
boston.mars <- earth(cmedv ~ ., data = boston, degree = 3)
pd <- partial(boston.mars, pred.var = c("lstat", "rm"))
plotPartial(pd)
plotPartial(pd, col.regions = colorRampPalette(c("red", "white", "blue")))
plotPartial(pd, contour = FALSE, zlab = "cmedv", drape = TRUE,
  colorkey = TRUE, screen = list(z = -20, x = -60))
```

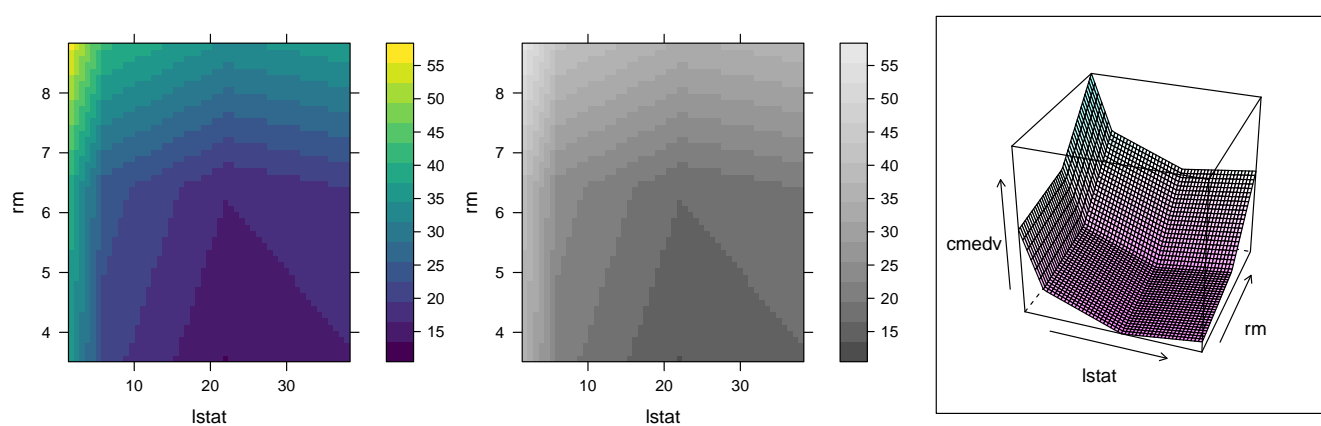
**Note:** the default color map for contour plots is the color blind-friendly `matplotlib` (Hunter, 2007) 'viridis' color map provided by the `viridis` package (Garnier, 2016).

### Avoiding extrapolation

It is not wise to draw conclusions from PDPs in regions outside the range of the training data. Here we describe two ways to mitigate the risk of extrapolation in PDPs: rug displays and convex hulls. Rug displays are one-dimensional plots added to the axes. Both `partial` and `plotPartial` have a `rug` option that, when set to `TRUE`, will display the deciles of the distribution (as well as the minimum and maximum values) for the predictors on the horizontal and vertical axes. The following snippet of code produces the left display in Figure 4.



**Figure 2:** Partial dependence of `cmedv` on `lstat` based on a random forest. *Left:* Default plot. *Right:* Customized plot obtained using the `plotPartial` function.

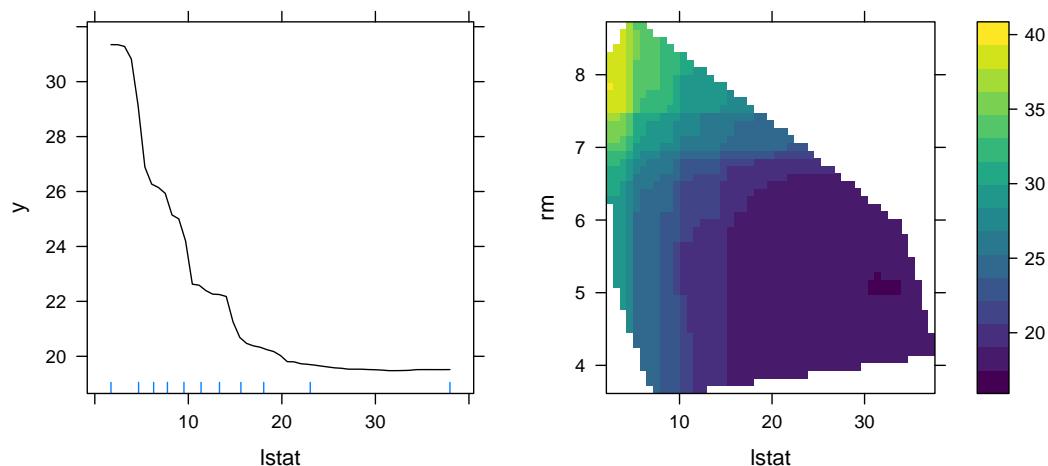


**Figure 3:** Partial dependence of `cmedv` on `lstat` and `rm` based on a third degree MARS model. *Left:* Default plot. *Middle:* Using a black and white color palette. *Right:* Using a 3-D surface.

```
partial(boston.rf, pred.var = "lstat", rug = TRUE, plot = TRUE)
```

In two or more dimensions, plotting the convex hull is more informative; it outlines the region of the predictor space that the model was trained on. When `chull = TRUE`, the convex hull of the first two dimensions of  $\mathbf{z}_s$  (i.e., the first two variables supplied to `pred.var`) is added to the plot; alternatively, you can set `chull = TRUE` in the call to `partial`, in which case only the region within the convex hull of the first two variables is plotted. Over interpreting the PDP outside of this region can be dangerous. The right display in Figure 4 was produced using:

```
partial(boston.rf, pred.var = c("lstat", "rm"), chull = TRUE, plot = TRUE)
```



**Figure 4:** Examples of PDPs with the addition of a rug display (left) and a convex hull (right).

### Addressing computational concerns

Constructing PDPs can be quite computationally expensive. Several strategies are available to ease the computational burden in larger problems. For example, there is no need to compute partial dependence of median home value using each unique value of `rm` in the training data (which would require  $k = 446$  passes over the data!). We could get very reasonable results using a reduced number of points. Current options are to use a grid of equally spaced values in the range of the variable of interest; the number of points can be controlled using the `grid.resolution` option in the call to `partial`. Alternatively, a specific set of values (e.g., quantiles of interest) can be supplied through the `pred.grid` argument. To illustrate, the following snippet of code computes the partial dependence of median home value on `rm` using each option; the results are displayed in Figure 5.

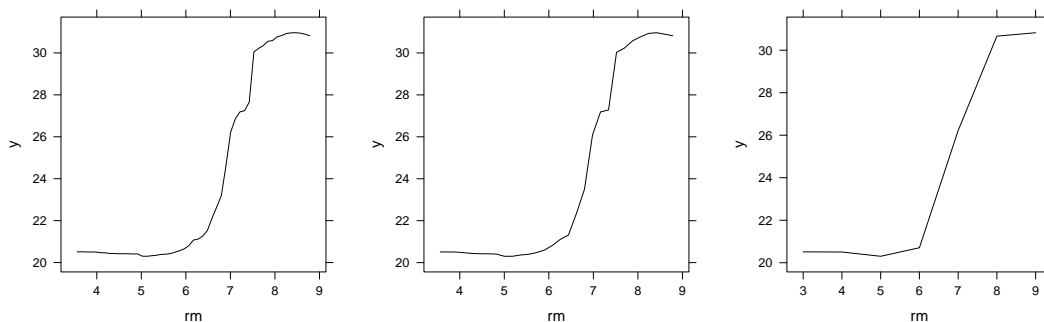
```
partial(boston.rf, "rm", plot = TRUE)
partial(boston.rf, "rm", grid.resolution = 30, plot = TRUE)
partial(boston.rf, "rm", pred.grid = data.frame(rm = 3:9), plot = TRUE)
```

The partial function relies on the `plyr` package (Wickham, 2011), rather than for loops. This makes it possible to request progress bars (e.g., `.progress = "text"`) or run `partial` in parallel. In fact, `partial` can use any of the parallel backends supported by the `foreach` package. To use this functionality, we must first load and register a supported parallel backend [e.g., `doMC` (Analytics and Weston, 2015a) or `doParallel` (Analytics and Weston, 2015b)].

To illustrate, we will use the Los Angeles ozone pollution data described in Breiman and Friedman (1985). The data contain daily measurements of ozone concentration (`ozone`) along with eight meteorological quantities for 330 days in the Los Angeles basin in 1976. The data are available from <http://statweb.stanford.edu/~tibs/ElemStatLearn/datasets/LAozone.data>. Details, including variable information, are available from <http://statweb.stanford.edu/~tibs/ElemStatLearn/datasets/LAozone.info>. The following code chunk loads the data into R:

```
ozone <- read.csv(paste0("http://statweb.stanford.edu/~tibs/ElemStatLearn/",
                        "datasets/LAozone.data"), header = TRUE)
```

Next, we use MARS to model ozone concentration as a nonlinear function of the eight meteorological variables plus day of the year; we allow for up to three-way interactions.



**Figure 5:** Partial dependence of cmedv on rm. *Left:* Default plot. *Middle:* Using a reduced grid size. *Right:* Using a user-specified grid.

```
ozone.earth <- earth(ozone ~ ., data = ozone, degree = 3)
summary(ozone.earth)
```

The model produced a generalized  $R^2$  of 0.79, similar to what was reported in [Breiman and Friedman \(1985\)](#). A single three-way interaction was found involving the predictors

- wind: wind speed (mph) at Los Angeles International Airport (LAX)
- temp: temperature ( $^{\circ}F$ ) at Sandburg Air Force Base
- dpgr: the pressure gradient (mm Hg) from LAX to Dagget, CA

To understand this interaction, we can use a partial plot. However, since the partial dependence between three continuous variables can be computationally expensive, we will run `partial` in parallel.

Setting up a parallel backend is rather straightforward. To illustrate, the following snippet of sets up the partial function to run in parallel on Unix-like systems<sup>1</sup> using the **doParallel** package.

```
library(doParallel) # load parallel backend
registerDoParallel(cores = 4) # use 4 cores
```

Now, to run `partial` in parallel, all we have to do is invoke the `.parallel = TRUE` option and the rest is taken care of by the internal call to **plyr** and the parallel backend we loaded. This is illustrated in following snippet of code which obtains the partial dependence of ozone on wind, temp, and dpgr in parallel. The result is displayed in Figure 6.

```
partial(ozone.earth, pred.var = c("wind", "temp", "dpgr"), plot = TRUE,
        chull = TRUE, .parallel = TRUE)
```

It is important to note that when using more than two predictor variables, `plotPartial` produces a trellis display. The first two variables given to `pred.var` are used for the horizontal and vertical axes, and additional variables define the panels. If the panel variables are continuous, then shingles<sup>2</sup> are produced first using the equal count algorithm (see, for example, `?lattice::equal.count`). Hence, it is probably better to use categorical variables to define the panels in higher dimensional displays when possible, as was done in the previous example.

## Classification problems

For classification problems, partial dependence functions are on a scale similar to the logit; see, for example, [Hastie et al. \(2009, pp. 369–370\)](#). Suppose the response is categorical with  $K$  levels, then for each class we compute

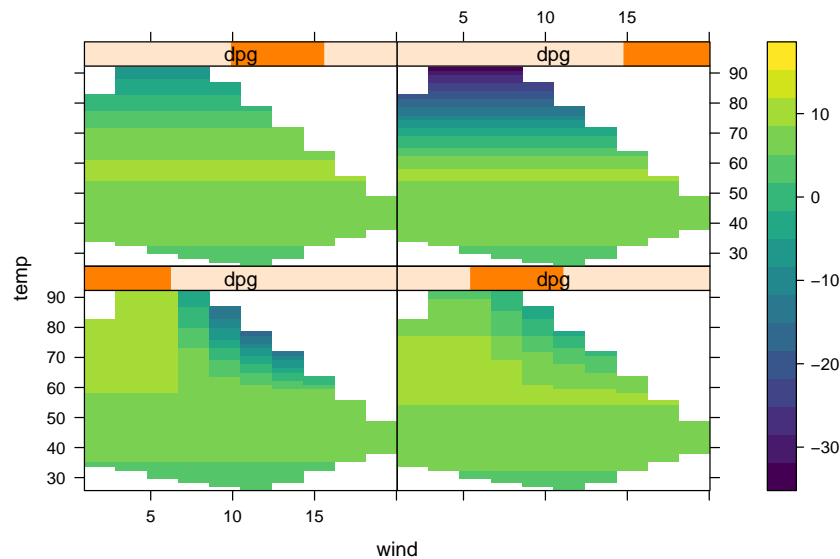
$$f_k(x) = \log[p_k(x)] - \frac{1}{K} \sum_{i=1}^K \log[p_i(x)], \quad k = 1, 2, \dots, K, \quad (3)$$

where  $p_k(x)$  is the predicted probability for the  $k$ -th class. Plotting  $f_k(x)$  helps us understand how the log-odds for the  $k$ -th class depends on different subsets of the predictor variables.

<sup>1</sup>This example will not run on Windows.

<sup>2</sup>A shingle is a special Trellis data structure that consists of a numeric vector along with intervals that define the "levels" of the shingle. The intervals may be allowed to overlap.





**Figure 6:** Partial dependence of ozone on wind, temp, and dpg. Since dpg is continuous, it is first converted to a shingle; in this case, four groups with 10% overlap.

To illustrate, consider Edgar Anderson’s iris data from the **datasets** package. The iris data frame contains the sepal length, sepal width, petal length, and petal width (in centimeters) for 50 flowers from each of three species of iris: setosa, versicolor, and virginica. We fit a support vector machine with a Gaussian radial basis function kernel to the data using the `svm` function in the **e1071** (the model parameters were determined using 5-fold cross-validation).

```
library(e1071)
iris.svm <- svm(Species ~ ., data = iris, kernel = "radial", gamma = 0.75,
               cost = 0.25, probability = TRUE)
```

**Note:** the partial function has to be able to extract the predicted probabilities for each class, so it is necessary to set the `probability = TRUE` option in the call to `svm`.

Next, we plot the partial dependence of Species on both Petal.Width and Petal.Length for each of the three classes. The result is displayed in Figure 7. Note that we had to supply the original training data to `partial` through the `train` option. This is necessary for fitted model objects that don’t store a copy of the training data (e.g., model fitting functions that don’t rely on a formula method).

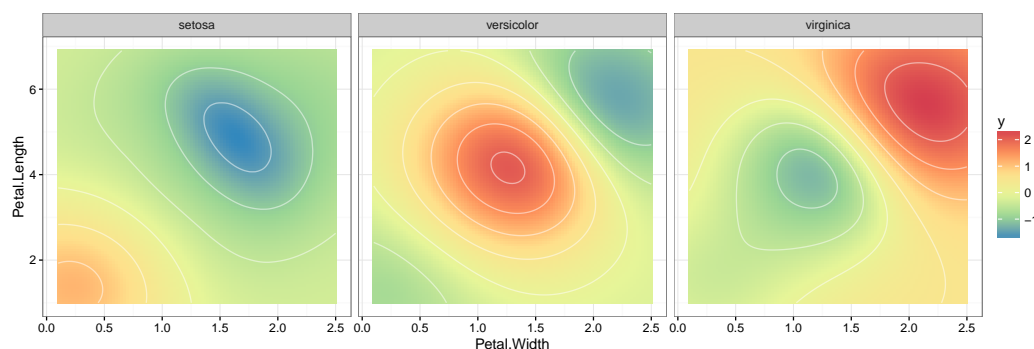
```
pd <- NULL
for (i in 1:3) {
  tmp <- partial(iris.svm, pred.var = c("Petal.Width", "Petal.Length"),
                which.class = i, grid.resolution = 101, .progress = "text")
  pd <- rbind(pd, cbind(tmp, Species = levels(iris$Species)[i]))
}

library(ggplot2)
ggplot(pd, aes(x = Petal.Width, y = Petal.Length, z = y, fill = y)) +
  geom_tile() +
  geom_contour(color = "white", alpha = 0.5) +
  scale_fill_distiller(palette = "Spectral") +
  theme_bw() +
  facet_grid(~ Species)
```

### Using partial with the XGBoost library

To round out our discussion, we provide one last example using a recently popular (and successful!) machine learning tool. XGBoost, short for eXtreme Gradient Boosting, is a popular library providing optimized distributed gradient boosting that is specifically designed to be highly efficient, flexible and portable. The associated R package **xgboost** has been used to win a number of [Kaggle competitions](#). It has been shown to be many times faster than the well known **gbm** package. However, unlike **gbm**,





**Figure 7:** Partial dependence of species on petal width and petal length for the iris data.

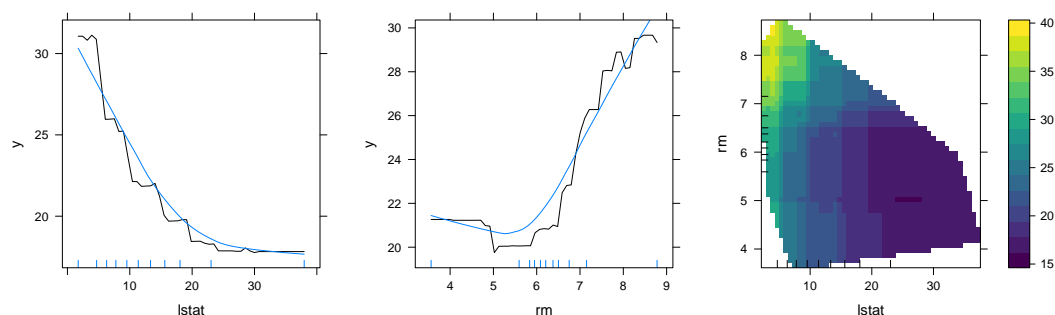
**xgboost** does not have built-in functions for displaying PDPs. Fortunately, the **pdp** package can be used to fill this gap.

For illustration, we return to the Boston housing data. The code chunk below fits an **xgboost** model to the boston data frame from Section 2.2; the tuning parameters were predetermined using 5-fold cross-validation. (Make sure you are using the latest version of **xgboost**: <https://github.com/dmlc/xgboost/tree/master/R-package>.)

```
set.seed(102) # for reproducibility
boston.xgb <- xgboost(data = data.matrix(subset(boston, select = -cmedv)),
  label = boston$cmedv, objective = "reg:linear",
  nrounds = 2000, max_depth = 3, eta = 0.01)
```

To use **partial** with the "xgb.Booster" class, we need to supply the original training data (minus the response!) in the call to **partial**. The following snippet of code computes the partial dependence of **cmedv** on both **rm** and **lstat**, individually and together. The results are displayed in Figure 8. **Note:** while **xgboost** requires the training data to be an object of class "matrix", "dgCMatrix", or "xgb.DMatrix", **partial** requires a "data.frame".

```
pred.list <- list("lstat", "rm", c("lstat", "rm"))
for (x in pred.list) {
  partial(boston.xgb, pred.var = x, plot = TRUE, smooth = TRUE, rug = TRUE,
    chull = TRUE, train = subset(boston, select = -cmedv))
}
```



**Figure 8:** PDPs for the top two most important variables in the Boston housing data using XGBoost. Compare this to the random forest results displayed in Figures 4-5.

## Summary

In this paper, we showed how to construct PDPs for various types of black box models in R using the **pdp** package. We also briefly discussed related approaches available in other R packages as well. Suggestions to avoid extrapolation and high execution times were discussed and demonstrated via examples. There is a bright future for **pdp** in terms of growth. For example, it would be useful to have the ability to construct PDPs for black box survival models—like conditional random forests with censored response.

## Acknowledgments

TBD.

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