

Statistical Learning Methods for Tree Classification using Remote Sensing Imagery

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Deeply grateful to...

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Abstract

The preface pretty much says it all.

Second paragraph of abstract starts here.

Introduction

0.1 Research Problem and Background

Western redcedar trees are evergreen trees that typically grow up to 75 feet tall and are located throughout the Pacific Northwest, making it an organism with a tolerance for shaded regions with moist environments. These trees are native to the land and have served many purposes to people and animals living in the vicinity of the trees, including medicinal, building, and habitat functions (Peterson, n.d.). Western redcedars were culturally important to indigenous peoples who valued the strong wood and utilized it for construction and everyday necessities (“Western red cedar,” n.d.). Over the past decade, reports of dead western redcedars have been increasing, suggesting that something other than natural causes is killing off this species. In general, western redcedars experience several hardships in surviving in the Pacific Northwest, with common causes of death such as forest fires, clearcutting, small animals eating the saplings, and harsh weather including strong winds that easily uproot the trees (Peterson, n.d.). Dying redcedar trees can be identified by their branches which turn brownish yellow or fall off completely. Another sign is that the top of a tree will turn brown and lose leaves (“Western redcedar Dieback,” n.d.). Losing this native tree would have a detrimental effect on animals in the area who rely heavily on the trees for their lifestyle. Scientists have speculated that western redcedar decline might be caused by recent dry summers, the spread of tree disease, insects, or other weather related events (“Western redcedar Dieback,” n.d.). Since this is a recent issue, there is not a lot of resources explaining the decline. This research aims to provide more insight into the cause of the western redcedar decline by first predicting the location of the western redcedar trees in Portland and then predicting their condition in terms of health. Having more insight helps to prevent further tree deaths and save the western redcedar species, which also extends to similar tree species and provides more knowledge about environmental changes in the Pacific Northwest region.

Modelling in this research will be conducted by combining information gathered from remotely sensed images with ground level information. There are several sources publishing research done using satellite imagery for land classification (Castelluccio, Poggi, Sansone, & Verdoliva, 2015) and for predicting tree species (Fricker et al., 2019), which will be the groundwork for this project’s application of remote sensing models to the specific topic of western redcedar mortality.

Remote sensing is a process for identifying the physical aspects of a large area of land by collecting and measuring the reflected light using airplanes or satellites. This method provides access to lot of information about the region of interest that would otherwise be limited from a human’s ground-level perspective. Remote sensing is used to detect characteristics on land as well as the sea. Some remote sensing land applications include tracking forest fires, volcanic eruptions, city growth, and also forest changes.

Satellites orbit the earth, carrying sensors that record levels of electromagnetic radiation detected by the reflection of the sun on the earth. The data collected is the measured radiation from different regions of the spectrum (visible light, infrared, ultraviolet, etc.). Depending on the surface, sunlight gets absorbed or reflected back

at the orbiting satellite. The remotely sensed data is available at certain resolutions depending on the instruments used. Spatial resolution is the area imaged by a satellite image, where a detailed image has smaller spatial resolution hence smaller pixels. Spectral resolution is a measure of the size of the wavelength interval, where smaller bands and smaller wavelengths improve the spectral resolution of an image and enhance the level of detail. Satellite images need to be corrected to combine the bands into an image depending on the analysis.

Imagery come in various band widths and different number of bands. Hyperspectral imagery has narrow bands of sizes 10-20nm and hundreds of bands. So, hyperspectral imagery is successful in distinguishing fine details, but at a cost of more complexity. Multispectral imagery has wider bands and up to 10 total bands. This type of imagery is not complex to work with, but misses some details that hyperspectral imagery detects.

Previous literature has been published in utilizing satellite imagery for predictive models, however, many issues arise in classifying land type through remote sensing mainly due to image quality. A single image can be composed of image strips taken over the course of multiple flight paths. Consequently, these images are taken at different times of day, which compiles into a single image with a lot of variation due to changes in the weather as well as the different angles of the sun's position (Castelluccio et al., 2015). Ideally, a solid classifying model surpasses any error introduced by imperfection in the satellite images. Common approaches to classification on satellite images include support vector machines, random forests, and decision trees. One study on land type classification explores the performance of convolutional neural networks (CNN) as classification models (Castelluccio et al., 2015).

Another related study identifies 7 tree species by applying a hyperspectral CNN model. This study was completed using field data with measured information about tree count, tree species, and mortality status and hyperspectral imagery data. The hyperspectral data is converted into the form of a tree canopy height model with circular polygons centered at individual tree canopies. The study analyzes the performance of CNNs for both hyperspectral imagery data and a Red-Green-Blue (RGB) subset of the hyperspectral imagery data. The results of the experiment conclude that training a CNN on the hyperspectral data outperforms the CNN trained on RGB data in classifying land type on the UC Merced Land Use dataset. The RGB model does not perform as well as the hyperspectral model in terms of distinguishing tree classes, but does perform around the same level of accuracy in terms of genus classification. For this project, the data comes from Planet.com and only has RGB and IR data available for the Portland region.

This thesis follows along with the methods in (Frick et al., 2019) for combining the satellite imagery data with ground data by creating spatial polygons and extracting pixel-level information to train classification models. The work in this thesis differs from the methods in the cited literature by using random forests and support vector machines as classification models instead of CNNs and uses the RGB model instead of hyperspectral model. Finally, this thesis applies established classification methods for satellite imagery data to answer the specific question about the cause of death of western redcedars in the Pacific Northwest. The models predict the loca-

tions for each tree species to facilitate the identification of any patterns in the species mortality over the past decade.

0.2 Overview

0.2.1 A Statistical Learning Approach to Identifying Location of Western Redcedars

This work combines RGB imagery data from Planet.com with ground level tree data from the RStudio `pdxTrees` library and applies random forest and support vector machine classification methods to predict the location of tree species (specifically western redcedars) in Portland, Oregon and model the condition of western redcedars to ultimately understand the cause of death in this species. Chapter 1 describes the two levels of data in this research: ground level and pixel level data, where the data comes from, and how the data is combined for modeling. Chapter 2 discusses the process behind training random forest and support vector machine models. Chapter 3 summarizes how the data is prepared for the final model and the outcome of the research followed by a discussion and conclusion.

Chapter 1

Data

There are two sources of data in this project. Data at a pixel level come from satellite images downloaded from Planet.com (<https://www.planet.com>), and data at the ground level come from library `pdxTrees` in RStudio (<https://github.com/mcconvil/pdxTrees>).

1.0.1 Imaging

Satellite images come from Planet.com's PS2.SD instrument found on Dove-R satellites that were launched in 2017. Satellites have an approximate frame size of 20 km x 12 km parallel to flight path. From a single pass, 4 to 5 overlapping framed scenes are combined to form a tile. To separate the light into red, blue, green, and near-infrared (NIR) channels, the telescope has a high-performance butcher-block filter made of 4 individual pass-band filters. Planet.com equates this pass-band filter with that of Sentinel-2.

As it orbits the Earth, the satellite captures continuous strips of single frame images that are split into a RGB frame and NIR frame. The butcher-block pass separately photographs the red, blue, green, and NIR bands and then combines all four to form an image. Images are downloaded already fully processed and ready to be analyzed. The process includes corrections for radiometric calibration, terrain distortions corrections, elevation corrections, and atmospheric corrections.

The images used in this research were taken on September 2nd, 2020 at an altitude of 475 km. Images are taken from the summer months to reduce the chance of rain clouds and capture peak greenness of the trees. The compiled multi-spectral image is composed of 4 bands: blue, green, red, and NIR, with center wavelengths 490 nm, 565 nm, 665 nm, and 865 nm respectively, with an average bandwidth of 41 nm. The average spatial extent is around 25 km by 8 km. The images cover the entire Portland area, with specific measurements of 26 km from west to east and 30 km from north to south. The spatial resolution of the pixels is approximately 3 meters.

Planet.com provides a way to download free images that are already processed and corrected for analysis. Some drawbacks to using this data are the limited number of downloads available per free account, the low spatial resolution, which decreases the performance of statistical models trying to predict tree species location, and the low

spectral resolution, so the satellite sensor cannot detect wavelengths in detail.

1.0.2 Ground Data

Ground level data is available from the ‘pdxTrees’ package in RStudio. This data was collected as part of the Portland’s Parks and Recreation’s Urban Forestry Tree Inventory Project, which collected park tree information from 2017 to 2019. Originally, the inventory project started with the goal of improving the city’s tree management plans. Under the guidance of US Forestry staff, volunteers around the city’s neighborhoods are trained in identifying tree species and provided with tools necessary to record tree measurements. Measurements in feet were made using diameter tape. The Portland park trees inventory consists of over 25,000 trees, each with information about tree location, size, species, and health. For the purposes of this project, the following variables were selected:

- Spatial information: Longitude and latitude of tree. Location of tree is recorded on a tablet with Collector for ArcGIS and recorders manually select the tree from a satellite image on the screen. The tree’s location is mapped over the satellite images in QGIS for analysis.
- Species: Tree genus, species, and common name. The tree species is used to train the model and predict the tree species of pixels in the test data.
- Crown size: Crown width from north to south, crown width from east to west, and the base height of the crown in feet. The crown of the tree is the tree’s above ground leaves, so the crown width is the longest horizontal distance that can be measured between the leaves of the tree. A measuring wheel is used to measure this distance.
- Tree Condition: Categorical variable with four categories (from best to worst), good, fair, poor, and dead. A tree was considered good if the tree is strong and has no apparent issues, fair if the tree is average condition with possibly a few dead branches, poor if the tree has major wounds and dead major canopy loss, and dead if the tree has no live leaves.

To prepare the ground level data for training the model, the data was filtered by species and crown size. The following tree species were identified and included: maples, oak, Douglas-fir, Western Redcedar, and sequoia. These species are all native to the Pacific Northwest, so they benefit the ecosystem and having a model of these trees will help track the species’ locations. Since the data is used to identify trees in the satellite images, the ground data was also filtered to only include the trees with a crown width from north to south of 20 feet or more, this ensures that there will be around 9 to 20 pixels per tree.

Portland’s Tree Inventory includes a variable for identifying trees including Western Redcedars as well as variables for filtering to include larger trees canopies (at least 6m in diameter) in the dataset, which helps construct a training set of tree polygons

over a satellite image. This data would be even better suited for the purposes of this project if the street trees subdata also contained variables for tree canopy size.

1.0.3 Training Data

Characteristics of Training Data

The training data at the ground-level was filtered by tree size and tree type. Trees with at least 20 feet in north to south crown width were included in the training dataset to ensure that the polygons around each tree contain around 9 to 20 pixels. Figure 1.1 plots the north to south crown width values against the west to east crown width values to present the ranges of crown widths in the entire pdxTrees parks dataset and to justify only using the north to south crown width for filtering the dataset since the two variables have a strong positive correlation.

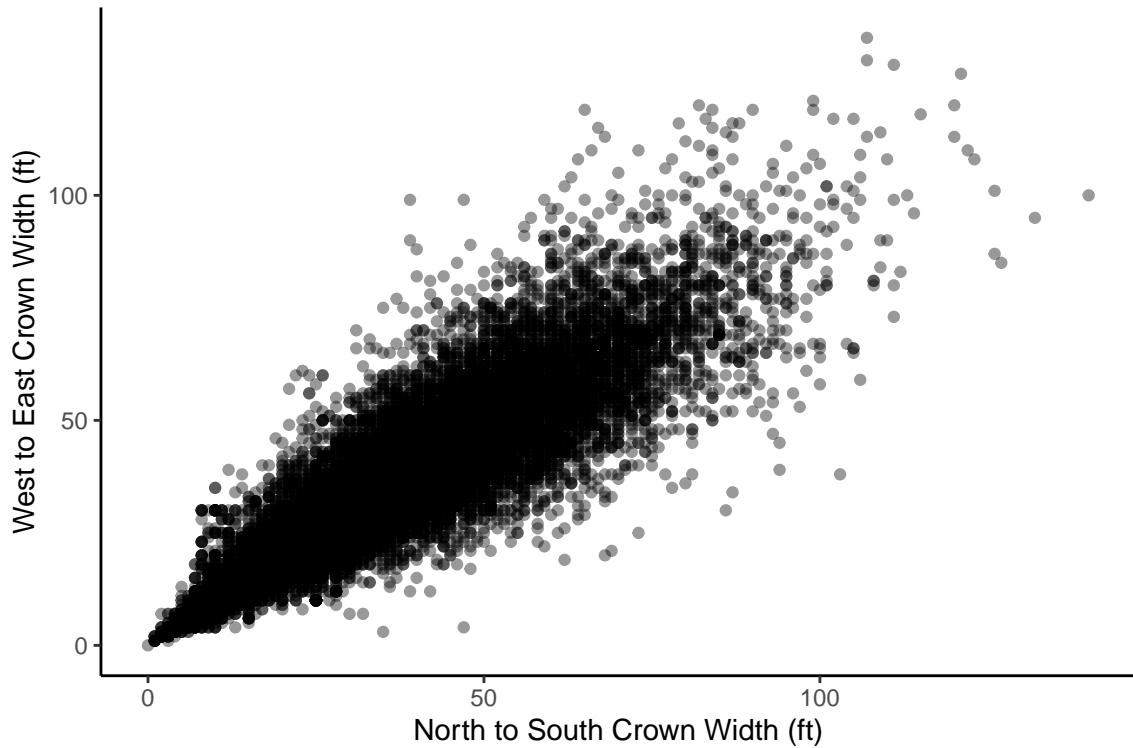


Figure 1.1: Plot comparing North to South Crown Width (ft) to West to East Crown Width (ft). A tree with a higher north-south crown width has a higher west-east crown width, so without loss of generality the north-south crown width variable is used to filter the dataset for larger trees only.

Six tree types were determined to have large enough canopies on average and appear frequently enough to construct a training dataset with plenty of observations for each species of tree: Douglas-Fir, English Oak, Giant Sequoia, Maple, Western Redcedar. Table 1.1 provides the number of trees under each common name category.

Table 1.1: Common tree names included in the tidy data and their total counts.

Tree Name	Tree Count
Douglas-Fir	6485
Norway Maple	1406
Western Redcedar	652
Bigleaf Maple	464
Giant Sequoia	315
English Oak	135

The largest group contains 6485 observations of Douglas-Fir trees, and the smallest group contains 135 observations of English Oak trees.

Creating Spatial Polygons

The training dataset was converted into shapefiles for each type of tree and exported into QGIS where polygons were manually drawn around 100 of the trees for each species. The point shapefile layers were displayed over the raster images downloaded from Planet.com in QGIS, which provided a guide point for locating individual trees. Then using QGIS's drawing tool, a polygon is carefully drawn around the tree canopy (see Figure 1.2 and Figure 1.3). Most polygons turned out to be four to six-sided polygons to retain the general shape of the tree canopy. Trees polygons were only created if the outline of the tree canopy was clearly visible or surrounded by other trees of the same species in order to avoid including pixels from the wrong species in that polygon. Also, the shadows of the trees are visible in the raster images, so the polygons were drawn with the intention of not including the tree shadow. For the five different species of trees, at least 100 polygons were drawn around trees of that type, with each polygon containing at least 6 pixels and at most 20 pixels. The polygon shapefiles were then exported into RStudio where the rest of the analysis was conducted.

Table 1.2 displays the number of polygons per raster strip and the amount of pixels from those polygons. Table 1.3 displays these counts in terms of tree species.

Table 1.2: Number of polygons per raster strip

Raster Strip	Polygons	n
a	197	3614
b	257	4887
c	36	359
d	8	98
e	177	2395



Figure 1.2: Polygons drawn around different tree species: maples (blue), sequoias (yellow), redcedars (pink). Note there are no polygons around trees with canopies that overlap with canopies of different tree species, like douglas-firs (green), to avoid misclassifying pixels from different tree species.



Figure 1.3: Polygons drawn around redcedar trees

Polygon Limitations

Ideally, the raster strips contain training trees of each species, and as a result, the polygons are evenly distributed across the raster strips, but due to limitations of our ground-level data as well as the coverage of the strips across the city of Portland, this is not achievable. Each of the raster strips contains some number of manually drawn polygons, however, raster strips ‘C’ and ‘D’ are missing some of the tree species. The raster strips do not all cover Portland to the same extent. Strips ‘A’ and ‘B’

Table 1.3: Number of polygons per tree type per raster strip

Raster Strip	Tree Name	Polygons	n
a	Bigleaf Maple	28	283
a	Douglas-Fir	35	432
a	English Oak	30	288
a	Giant Sequoia	26	336
a	grass	1	1158
a	Norway Maple	48	747
a	Western Redcedar	29	370
b	Bigleaf Maple	26	279
b	Douglas-Fir	38	604
b	English Oak	62	618
b	Giant Sequoia	32	399
b	grass	3	1579
b	Norway Maple	70	1026
b	Western Redcedar	26	382
c	Bigleaf Maple	9	55
c	Douglas-Fir	19	202
c	English Oak	2	39
c	Western Redcedar	6	63
d	Douglas-Fir	5	56
d	Western Redcedar	3	42
e	Bigleaf Maple	37	268
e	Douglas-Fir	6	86
e	English Oak	6	80
e	Giant Sequoia	54	701
e	grass	1	186
e	Norway Maple	37	567
e	Western Redcedar	36	507

Table 1.4: Variables and pixel values included in the pixels dataset

ID	red	green	blue	ir	rstrip	id_type	Genus	Species	Cmnn_Nm	Cr_W_NS	Cr_W_EW	Crw_B_H	Condtn	St_Wdth	ndvi
43	4707	4125	3187	6627	a	acpl	Acer	ACPL	Norway Maple	28	28	8	Fair	NA	0.1694018
43	4384	3860	2864	7111	a	acpl	Acer	ACPL	Norway Maple	28	28	8	Fair	NA	0.2372336
43	4446	3803	2807	7443	a	acpl	Acer	ACPL	Norway Maple	28	28	8	Fair	NA	0.2520818
43	4619	3911	2859	7801	a	acpl	Acer	ACPL	Norway Maple	28	28	8	Fair	NA	0.2561997
43	4695	4060	3201	5996	a	acpl	Acer	ACPL	Norway Maple	28	28	8	Fair	NA	0.1216911
43	4344	3719	2728	6747	a	acpl	Acer	ACPL	Norway Maple	28	28	8	Fair	NA	0.2166622

encompass most of the city while strips ‘C’ and ‘D’ only make up a small portion that covers the corners of the city not reached by strips ‘A’ and ‘B’.

The polygons are drawn with the intention of outlining pixels for a known tree species so that the extracted pixels truly represent that species and so that there is a sufficient amount of pixels per species, however, limitations of the satellite images’ resolution make some error inevitable. Many parks trees have canopies that overlap, which cause some polygons to contain pixels from trees of different species. Other pixel errors might come from including pixels that are cast in shadow by the polygon’s training tree or by surrounding infrastructure. To avoid these errors, polygons with uncertain canopies are cross-checked with the satellite filter on Google Maps. For example, a point in QGIS that is indicated as a douglas-fir tree might appear as one tree on the low-resolution image from Planet.com, but a closer look at Google Maps shows that it is actually two trees in close proximity. Having low spatial resolution decreases the number of pixels available within a polygon, but that has to be balanced out with the need for a large number of pixels. With regards to the number of pixels per tree species, having around the same number of tree polygons per species results in differing amounts of pixels per species due to the different average sizes of tree canopies for different species. After drawing the images, the pixels totals per tree species is computed to ensure that the training pixels data contains at least 800 pixels per tree species.

Combining Ground-level Data with Pixel-level Data

The first spatial join in RStudio was conducted to match up each Spatial Polygon with a point in the training dataset. Then the raster images were loaded and joined with the polygons to extract the pixel values inside each polygon for all 4 bands (red, green, blue, infrared). Ultimately this turned into a pixel table with rows representing each pixel with its corresponding polygon, light reflection intensity values for all 4 bands, and the ground information about that tree. Table 1.4 displays the first few entries of the pixels dataset. Table 1.5 contains the summary of the total number of pixels for each tree type. Figure 1.4 displays the range of reflection intensity pixel values per tree type. Each tree species has similar ranges of values over the red, green, blue, and infrared bands. Of the species included in the training data, the Giant Sequoia trees appear to have higher density counts for the red, green, and blue bands.

Table 1.5: Number of pixels per tree type

Tree Name	Pixels Count
Bigleaf Maple	885
Douglas-Fir	1380
English Oak	1025
Giant Sequoia	1436
grass	2923
Norway Maple	2340
Western Redcedar	1364

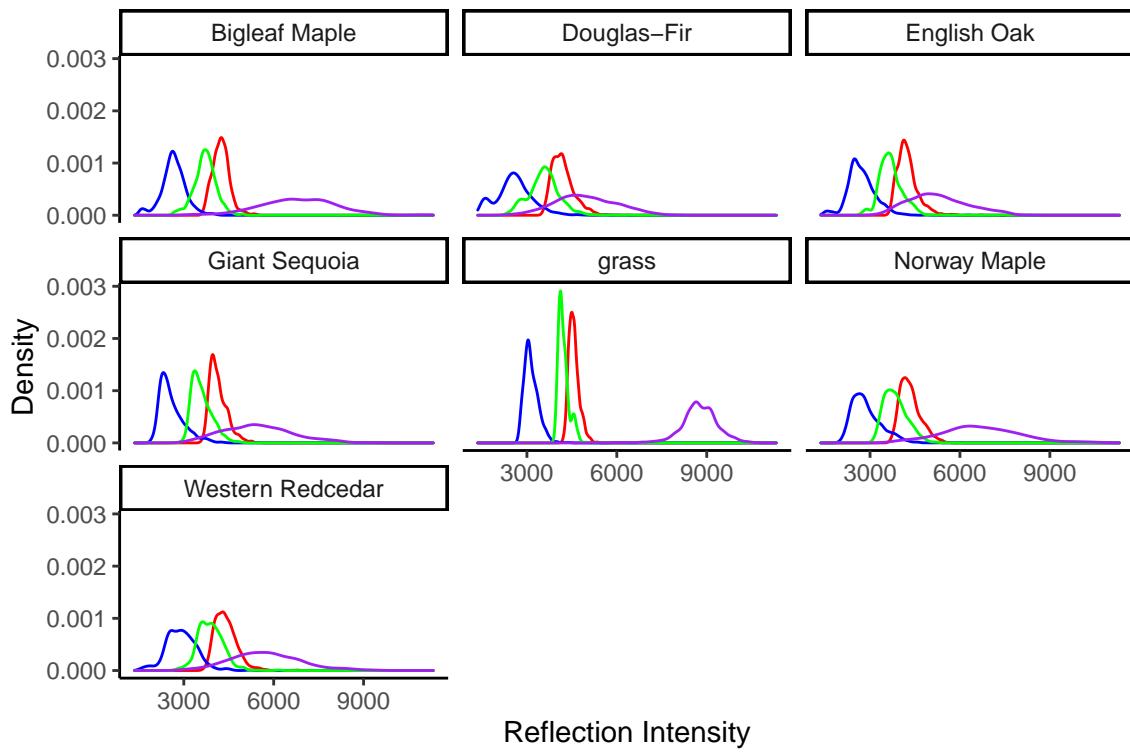


Figure 1.4

1.1 Preparing Raster Images

The training model has to be applied to the entire raster image to predict the location of Western redcedars, however, extracting all the pixels from raster images is a slow process. Filtering the rasters to keep vegetation and to recognize the difference between grass pixels and tree pixels alleviates the computational intensity of extracting all the pixels and improves the performance of the models. To reduce the size of computations, all five raster strips are masked and cropped to only include pixels within the city boundary of Portland, since that is the region of interest, and that is the extent of the ground level data.

1.1.1 Masking Vegetation

The satellite images in RStudio need to be masked to reduce the number of pixels that the model has to classify and prevent the chance of a non-vegetation surface being predicted as a tree. A common mask applied to raster images is a Normalized Difference Vegetation Index (NDVI) mask. This index is a measure of the greenness of a pixel, with higher values indicating vegetation and lower values indicating infertile areas such as a rock. The formula for NDVI is $NDVI = \frac{NIR - Red}{NIR + Red}$. Figure 1.5 displays the density of NDVI values over the raster images. To remove the pixels that are not vegetation, an NDVI threshold is determined to be 0.00 to create a mask dividing the pixels into vegetation (1) and non-vegetation (0). When the trained model is applied to the raster image, it is applied to the masked raster image that only keeps pixels with NDVI mask values of 1.

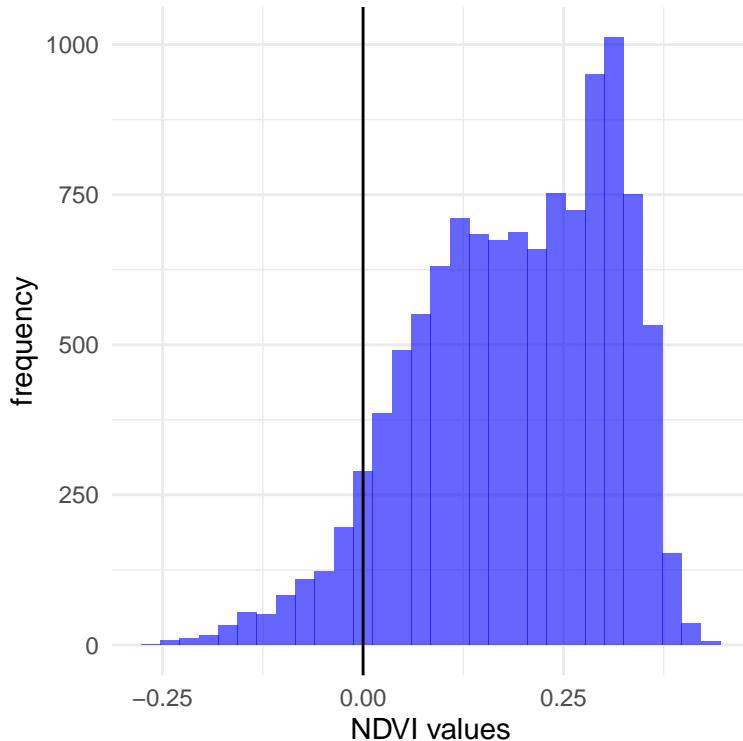


Figure 1.5: Normalized Difference Vegetation Index (NDVI) is a measure of the greenness of a pixel. The histogram displays the frequency of the tree pixels' NDVI values. A good threshold for masking out the nonvegetation pixels is an NDVI of 0.

1.1.2 Masking Grass

A second mask is applied to the NDVI masked raster image to distinguish tree pixels from grass pixels. Inspired by previous research, a grass index was created by averaging the values of the four bands, $GRASS = \frac{RED + BLUE + GREEN + NIR}{4}$ (Qian, Zhou, Nyctch, Han, & Li, 2020). Grass polygons were created in QGIS to add a grass at-

tribute in the pixels dataset. Five fields of grass were outlined as polygons in QGIS. In RStudio, the pixels were extracted from the grass polygons for a total of 2,923 grass pixels. Figure 1.6 displays how the average band values for the grass pixels differ from the other tree types average band values. Based on the figure, pixels in the raster image that have an average band value above 4900 are filtered out. Masking the grass pixels ensures that a field of grass will not be classified as a tree when applied to the entire raster image

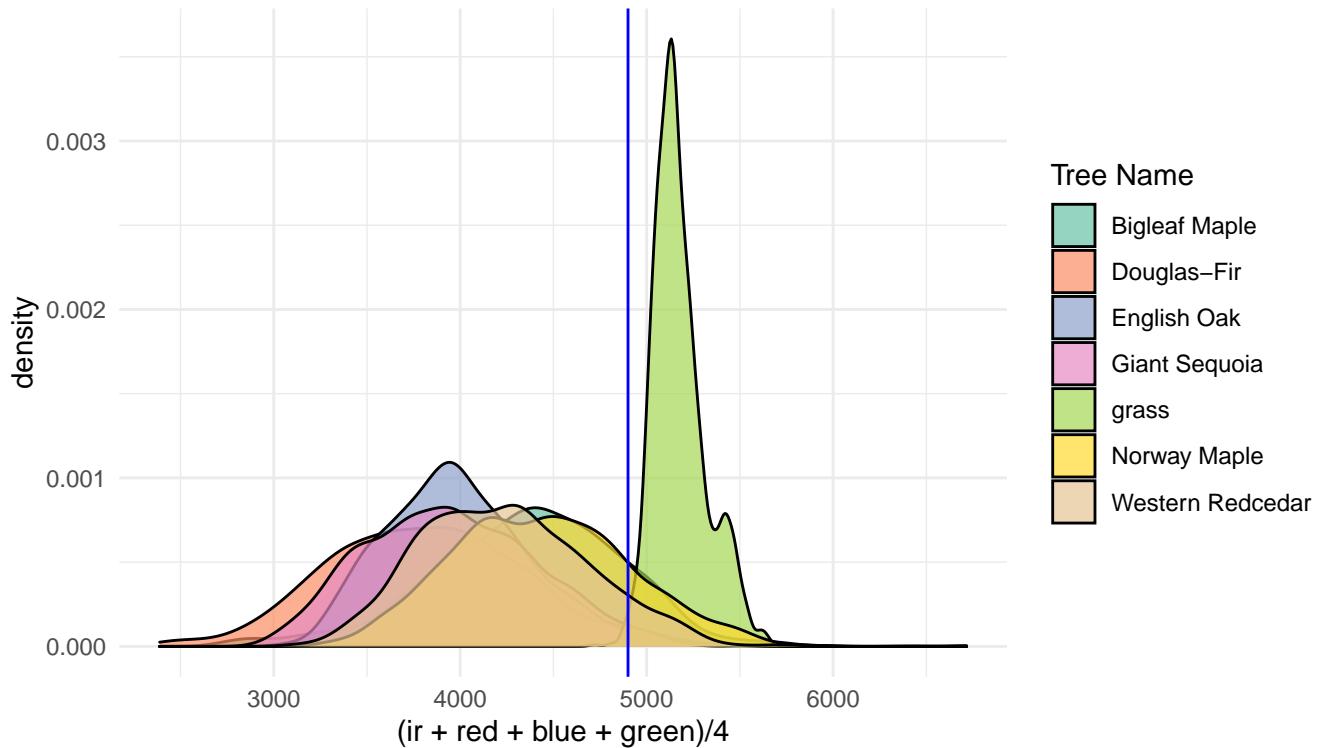


Figure 1.6: Average of band values distribution across different tree types in comparison with grass pixel average band values. Most grass pixels have an average value above 4900, so a mask is applied to the raster image to remove pixels with average values above this threshold.

Figure 1.7 displays a raster strip with every pixel included (top), only the pixels with NDVI values above 0.00 (middle), and only pixels with NDVI values above 0.00 and average band values below 4900 (bottom). Ideally, the bottom image is left with only the tree pixels in the image.

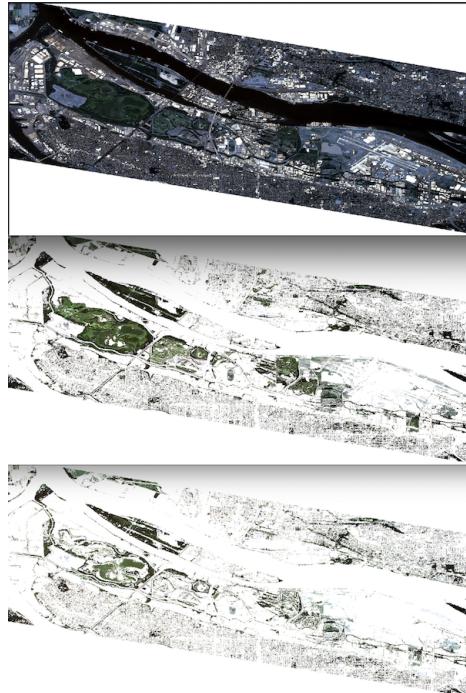


Figure 1.7: A raster strip with every pixel included (top), only the pixels with NDVI values above 0.00 (middle), and only pixels with NDVI values above 0.00 and average band values below 4900 (bottom)

1.1.3 Masking Limitations

The low resolution of the raster images cause shapes to appear blurry, and especially tree canopies that are close together appear as a single formation.

Masked raster images were compared in QGIS and, in general, the masks are successful in keeping the forest pixels while still removing building structures and roads. If a field of grass is a brownish color, it gets removed, but some greener fields are not masked out. The masks consistently remove river pixels but not lake pixels if the lake is a greenish color. For example, Smith Lake in North Portland is considered an urban wetland, which signifies a body of water with a lot of vegetation. From the perspective of a satellite, this looks like a field of grass. However, the average band pixel values are less than 4900, so they are not masked out by the final GRASS mask stage, and instead are treated as tree pixels in the model (see Figure 1.8).

To address leftover grass pixels in the filtered raster image, access to lidar data would provide a method of determining the height of certain pixels. Setting a height threshold would ensure that grass pixels are removed while higher vegetation pixels are kept. This would also address the masks' limitations in removing lakes with high vegetation, since those pixels would not meet the threshold for height.

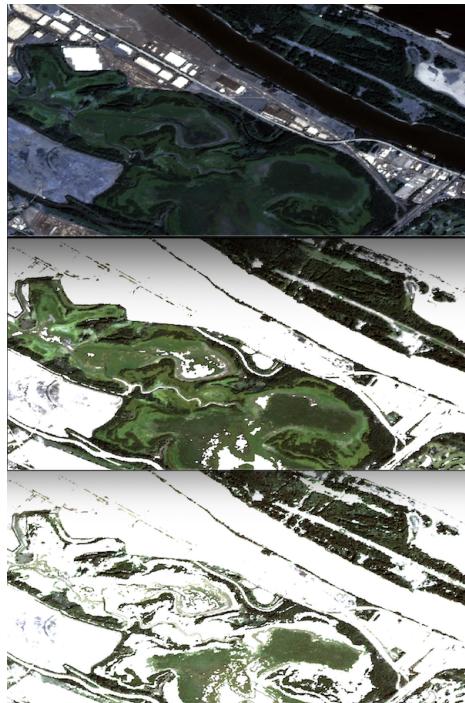


Figure 1.8: Satellite view of Smith Lake (top), after NDVI mask applied (middle), and NDVI mask and GRASS mask applied (bottom). The final masked image still contains parts of the lake because the aquatic vegetation fits the pixel values descriptions set by the masks.

Chapter 2

Methods

2.1 Current Methods

Since the Western redcedar decline is recent, there are currently no published work for identifying tree species like Western redcedars for the Pacific Northwest using satellite imaging. However, a similar study identifies tree species for a strip of land in California by combining ground-level data with satellite images (Fricker et al., 2019). Their methods involved drawing polygons around rastered images that are layered with ground data as explained in Chapter 1, but the pixels from the polygons are used to train a convolutional neural network (CNN) model instead of the random forests and support vector machines in this study. The CNN model was then evaluated with k-fold cross validation. CNN models are appropriate in a classification setting and when working with satellite images because they account for spatial relations, which is likely to appear in classifying tree species. This study follows the methods of preparing the data for modelling as well as the method of cross validation to evaluate the performance of the random forests and support vector machines models. Random forests are used in a classification setting or regression setting and have the advantage of being simpler to train than CNN models and still performs at a similar level. Random forests use random feature selections to create decision trees and increase the number of correctly classified observations. Since this research strives to provide a computationally simple method of estimating tree locations to track changes in the ecosystem, a random forest is suitable for the study. If a researcher was interested in making more precise predictions, a CNN would require more time but would produce better estimates.

2.1.1 Random Forest

In building a decision tree for classification, recursive binary splits are made by minimizing the Gini index (G), the total variance across K classes:

$$G = \sum_{k=1}^K \hat{p}_{mk}(1 - \hat{p}_{mk}),$$

where \hat{p}_{mk} is the proportion of observations in the m^{th} region from the k^{th} class (James, Witten, Hastie, & Tibshirani, 2013). A small Gini index value represents a region containing mostly observations from a single class, while a large Gini index value represents a lot of variation across classes. Then the decision tree is constructed by repeatedly considering different attributes and making splits where the Gini index is minimized. Note that this top-down, greedy approach tends to overfit the training data, so finding the decision tree that performs the best on test data involves a cost-complexity pruning algorithm to obtain smaller trees and apply k-fold cross validation to choose the best tree that minimizes the average error. K-fold cross validation is performed when training the models by splitting the data into $k = 10$ folds and training ten models each time withholding one of the ten folds until the process goes through every fold. This process aids in understanding how the models will perform against data that has not been included in training the model.

In the classification setting, a stronger predictive model than a decision tree is a random forest, which uses an element of randomness to decorrelate bootstrapped decision trees. Bootstrapped samples from the training dataset are used to construct decision trees, where splits are based on minimizing the Gini index when selecting from a random sample m of attributes from all p available. In common random forest applications, the number of attributes considered at a split is $m \approx \sqrt{p}$. Theoretically, by forcing only a subset of attributes to be featured in the tree, this expands the number of possible subtrees that might not have been achieved by the top-down, greedy approach. Hence, random forests provide a method of predicting classes with low variance while also keeping bias at a minimum.

2.1.2 Support Vector Machine

Another method investigated in this study imagines the multidimensional data in space and classifies the data using hyperplanes positioned to minimize misclassification error and maximize distance from observations to separate the data (James et al., 2013). Support vectors are vectors in space that represent the subset of observations which influence the hyperplane classifier. For example, in two dimensions, imagine a line separating A's one one side and B's on the other side. This line is the maximal margin hyperplane if it is the farthest possible from the observations and also separates the two classes perfectly. Observations that are closer to the hyperplane (support vectors) have more influence over the line if they move than observations that are further. Consequently, more support vectors indicates a classifier with lower variance and high bias. A tuning parameter can be altered to expand or shrink the margin of error surrounding the hyperplane (how far away the observations have to be from the plane) and decide the level of tolerance for misclassifying observations.

Support vector machines can be applied in a linear or nonlinear setting with the use of kernels. A kernel is a function $K(x_i, x_{i'})$ of two observations specifying the similarities between two points. Given data with n observations, let x_i be an observation, where $i = 1, \dots, n$. A linear kernel is

$$K(x_i, x_{i'}) = \sum_{j=1}^p x_{ij} x_{i'j}$$

with a support vector classifier

$$f(x) = \beta_0 + \sum_{i=1}^n \alpha_i \langle x, x_i \rangle.$$

The parameters α_i and β_0 are estimated using the inner products of the observations, but $\alpha_i \neq 0$ if and only if x_i is a support vector. Note that this function is similar to a linear regression function. Extending this model to a nonlinear context involves a similar approach as adding nonlinear terms to a regression model.

The polynomial kernel of degree d is defined to be

$$K(x_i, x_{i'}) = \left(1 + \sum_{j=1}^p x_{ij} x_{i'j} \right)^d$$

with support vector classifier

$$f(x) = \beta_0 + \sum_{i=1}^n \alpha_i K(x, x_i).$$

For some positive constant k , the radial kernel is defined to be

$$K(x_i, x_{i'}) = \exp \left(-\gamma \sum_{j=1}^p (x_{ij} - x_{i'j})^2 \right),$$

which makes the support vector machine more sensitive to nearby observations as opposed to further observations.

Support vector machines are less sensitive to outliers and successful predictors of categorical variables in high-dimensional data. A support vector machine does not perform well for data with observations that overlap and if the kernel is incorrectly selected.

2.2 Training Models

To model the tree type of the pixels in raster image, a random forest model and a support vector machine model are trained using the pixels training data. First the data is separated into a training set and then a test set by a ratio of 70% training data and 30% testing data. The `dplyr` function `slice_sample` was used to take a stratified sample from the entire pixels dataset where 70% of each tree species in an individual raster strip was randomly selected. For example, the training set consisted of 70% of the bigleaf maple trees in raster strip ‘A’, 70% of bigleaf maple trees in raster strip ‘B’, and so on. From the 70% training data, 500 observations for each tree species are extracted to obtain an even distribution of species.

The selected variables for training the models are the four bands (red, green, blue, infrared) and an NDVI variable, and the predictive variable is the tree species name. Some predictor variables were manually computed using ratios of bands. Typically, a long over short band ratio provides some more information to train the model. The following bands were created: red/blue, ir/red, red/green, and blue/green. Examining a correlation matrix with these predictors reveals possibly useful and not too highly correlated variables are red/green, red/blue, and blue/green (see Figure 2.1).

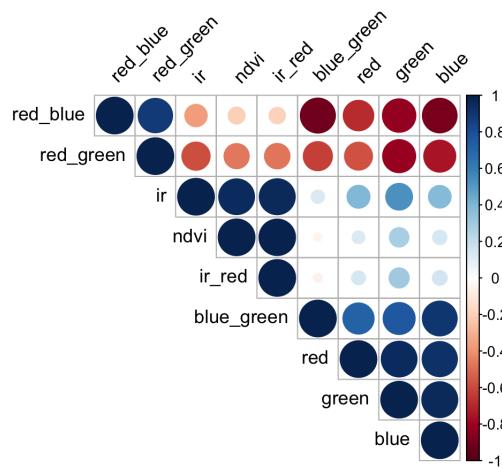


Figure 2.1: Correlation matrix of possible predictor variables to include in the model. Size and shade intensity represents stronger correlation (red/blue and blue are highly negatively correlated, while NDVI and green have almost no correlation). Positive correlation indicates higher values of one variable results in higher values of the other variable, while negative correlation indicates higher values of one variable results in lower values of the other variable.

To address the small number of predictor variables available to predict the large number of classes (7 tree species), some species are grouped together. The Bigleaf Maple, Norway Maple, and English Oak species are grouped under a new “Broadleaf” tree category. This reduces the number of classes to 5: Broadleaf, Douglas-Fir, Giant Sequoia, Grass, and Western Redcedar. Both the full 7 tree species training data and the grouped (5 class) species training data are used to train the models.

The `caret` package (Kuhn, 2020) in RStudio stands for classification and regression training, and it contains functions for facilitating the process of creating training models in R and determining the tuning parameters.

2.2.1 Random Forest Tuning

The `train` function from the `caret` package (Kuhn, 2020) with the specified `method = ranger` option is used to train a random forest model on the training dataset to predict the tree species. The function takes parameters for data, the predictive variable, and the method of training the model. To perform k-fold cross validation, the `trainControl` function is used to create an object that tells the `train` function to perform cross validation 10 times with a different fold left out each time. The function automatically tries multiple number of tries at each split (`mtry`) and determines the optimal number through cross-validation. Training preliminary models with varying number of predictors provided a sense of the number of predictors to include in the training data that still add information to the model but avoid adding redundancy or multicollinearity to the model. For example, just having 5 predictors in the model results in a random forest that only selects one variable for each split, while 12 predictors guarantees an issue of multicollinearity. Models were considered with total predictors around 7 or 8 predicting classes of size 7 (individual species) or 5 (oaks and maples grouped together):

- 7 predictor model: red, green, blue, infrared, NDVI, $\frac{\text{red}}{\text{green}}$, $\frac{\text{blue}}{\text{green}}$
- 8 predictor model: red, green, blue, infrared, NDVI, $\frac{\text{red}}{\text{green}}$, $\frac{\text{red}}{\text{blue}}$, $\frac{\text{blue}}{\text{green}}$

7 predictor Random Forest

The 7 predictor random forest model uses the following predictors: red, green, blue, infrared, NDVI, $\frac{\text{red}}{\text{green}}$ and $\frac{\text{blue}}{\text{green}}$ band values predictors. Figure 2.2 displays the accuracy of the model predicting all 7 tree classes and the grouped 5 classes training set with a range of `mtry` values. The 7 predictor model with the highest accuracy (determined by cross-validation) randomly selected 1 predictor at each split with an accuracy of 0.44. The variable with the highest predictive power was infrared band values with NDVI values with the second highest predictive power. For the grouped data, the optimal `mtry` value of 6 produced an accuracy of 0.57. The most important variable for this model was infrared followed by NDVI values.

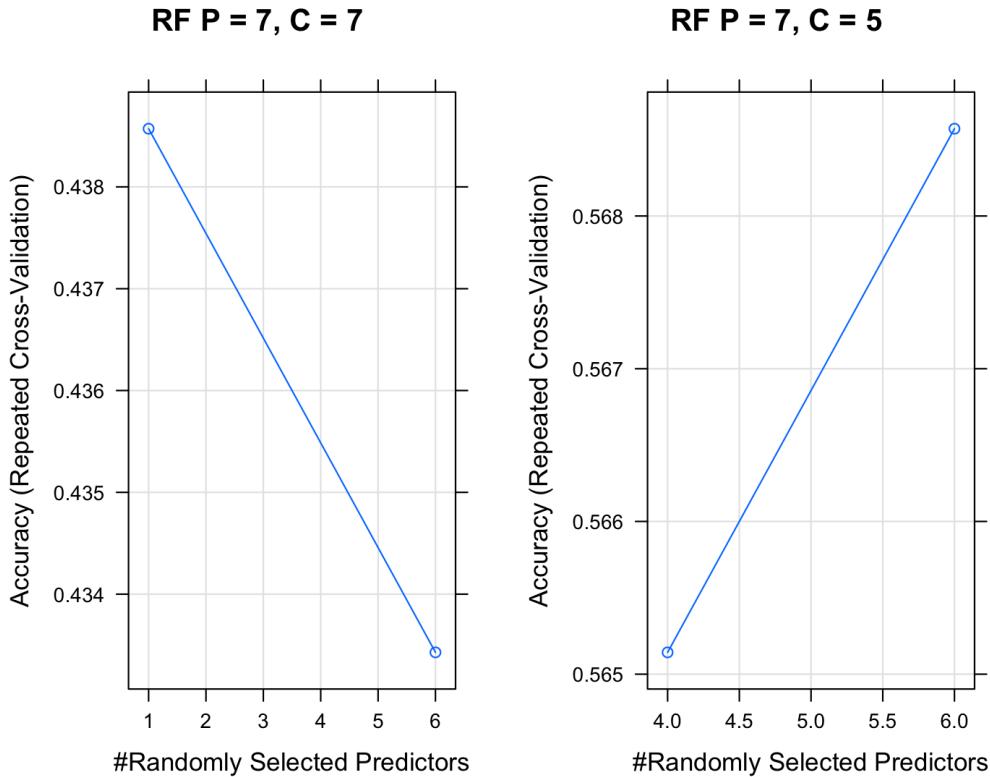


Figure 2.2: Accuracy of the 7 predictor trained random forest model for different mtry numbers over 10-fold cross-validation on 7 class and 5 class training data. (P = Predictors, C = Classes)

8 predictor Random Forest

The 8 predictor random forest model added a ^{blue} _{green} predictor to the 7 predictor model. Figure 2.3 displays the results of a different number of randomly selected predictors to be considered at each split in the tree for 7 class data and 5 class. For modelling all 7 classes, the optimal mtry value was 2 with an accuracy of 0.44. The variables with the higher predictive power in decreasing order were infrared, NDVI, and ^{red} _{green} band values predictors. The model using grouped data with 4 tree classes, had an optimal mtry value of 4 with an accuracy of 0.57. The most important variable was infrared.

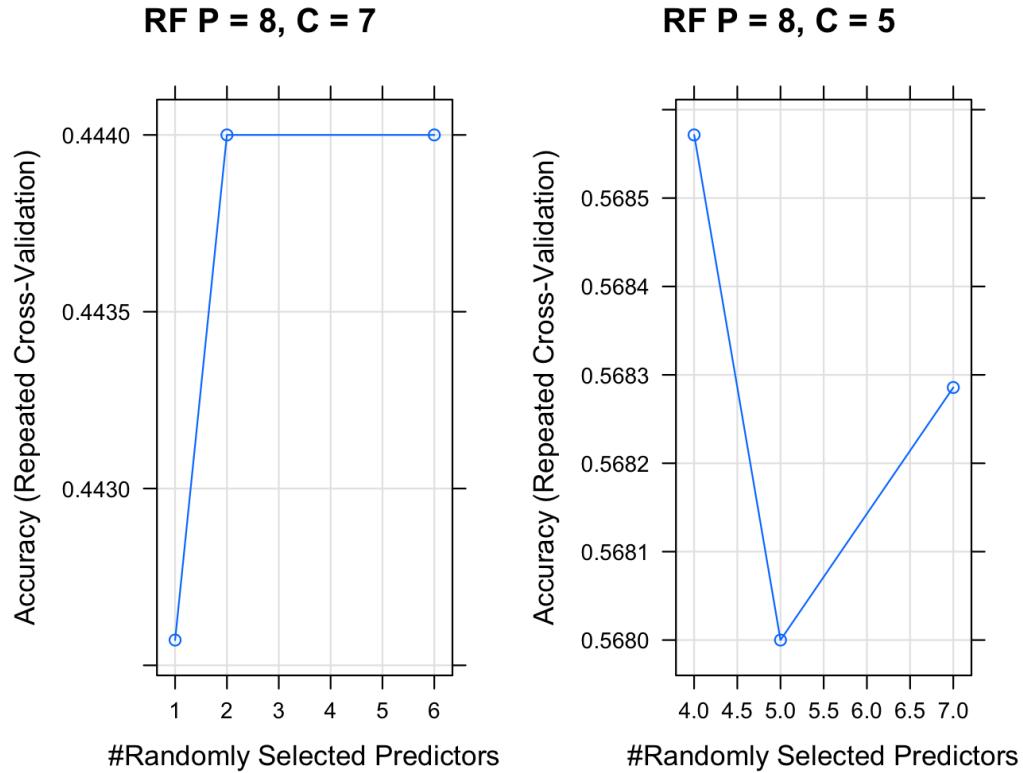


Figure 2.3: Accuracy of the 8 predictor trained random forest model for different mtry numbers over 10-fold cross-validation on 7 class and 5 class training data. (P = Predictors, C = Classes)

Best Random Forest

Figure 2.4 compares the results of the models based on overall model accuracy in predicting the test data.

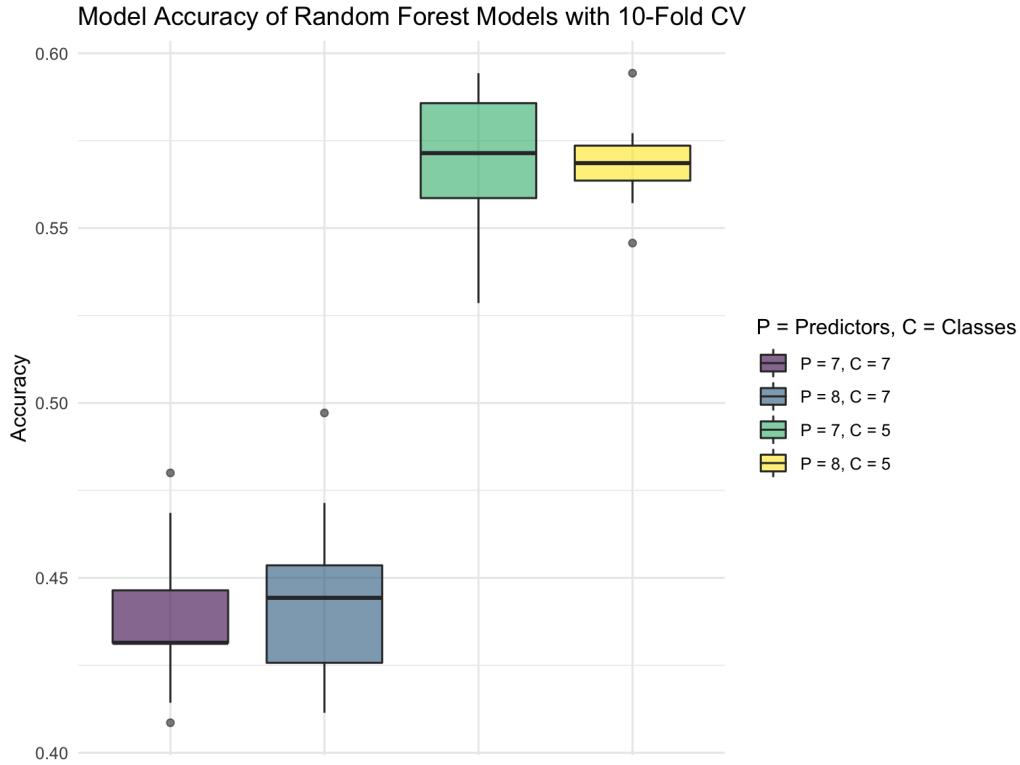


Figure 2.4: Comparison of results from random forest models with different predictors included

For modelling the 7 individual tree classes, the model with the 8 predictors included has a slightly higher maximum accuracy overall than the 7 predictor model. A confusion matrix located in the appendix was constructed for each model averaging the entry counts over all ten cross validation resamples to investigate model performance just predicting Western redcedar tree pixels. Both the 7 and the 8 predictor models performed similarly classifying Western redcedar pixels, the 7 predictor model had a prediction accuracy of 32%, the 8 predictor model had a prediction accuracy of 33%. Out of both models, Western redcedars tend to be inaccurately classified under douglas-firs, english oak, and norway maple trees.

Modelling only 5 tree classes with the grouped data produced higher accuracies than the full classes models. The 7 predictor model had a larger variation than the 8 predictor model, but the 8 predictor model produced the highest accuracies of 0.57 through cross-validation. The 7 predictor model accurately predicted Western redcedars 37% percent of the time, while the 8 predictor model predicted redcedars with an accuracy of 36%. Most inaccurately predicted redcedars were classified under the broadleaf category. Models on both training sets predicted grass pixels with the highest accuracy.

Table 2.1 compares the hyperparameters accross all random forest models.

Table 2.1: Accuracy of random forest models with different number of predictors and mtry determined by ten-fold cross validation.

Random Forest Model	Accuracy	Kappa	mtry
P = 7, C = 7	0.4386	0.3450	1
P = 8, C = 7	0.4440	0.3513	2
P = 7, C = 5	0.5686	0.3803	6
P = 8, C = 5	0.5686	0.3798	4

2.2.2 Support Vector Machine Tuning

The `train` function has the `method = svmLinear/svmRadial/svmPoly` option to train a SVM model on the dataset to classify pixels into species of trees. For multiple class SVM training on k classes, this function classifies “one-against-one” by training $k(k - 1)/2$ binary classifiers. This option also has parameters for cost, loss function, class weights, and normalized variables. For this project, three support vector machines were trained: linear, radial basis, and polynomial basis, and the preprocess setting for all three models were specified to center and scale (in R: `preProcess = c("center", "scale")`), which standardizes the variables. This is commonly performed with SVM models because the scale of the variables influences the optimal hyperplane decision. Since the light intensities extracted from the band data are only positive values, it is reasonable to consider normalizing the variables by setting the maximum value in the variable to 1 and the minimum value to 0, however, in comparing the SVM models under this preprocess setting (in R: `preProcess = "range"`) to the standardized setting, the accuracy of the models is not as high as the accuracy of the standardized variable models. The `caret` package selects the best cost tuning parameters based on accuracy through cross-validation.

Linear Support Vector Machine

The linear kernel SVM model had the highest training accuracy at 0.41 when cost parameter held constant at a value of 1. The final model contained 3020 support vectors. The linear kernel SVM model on the 5 classes data held the cost parameter constant at 1 predicted the grouped tree classes with an accuracy of 0.55 contained 2742 support vectors.

Radial Support Vector Machine

Figure 2.5 displays the results of different cost parameters on training accuracy for the radial basis kernel function. The radial kernel SVM model on the 7 classes data set performed best when cost was 32 and $\sigma = 0.23$ for a training accuracy of 0.46. The final model had 2889 support vectors. The final values for the radial model on the 5 classes data set were $\sigma = 0.24$ and cost = 1 with an accuracy of 0.59 and 2612 support vectors.

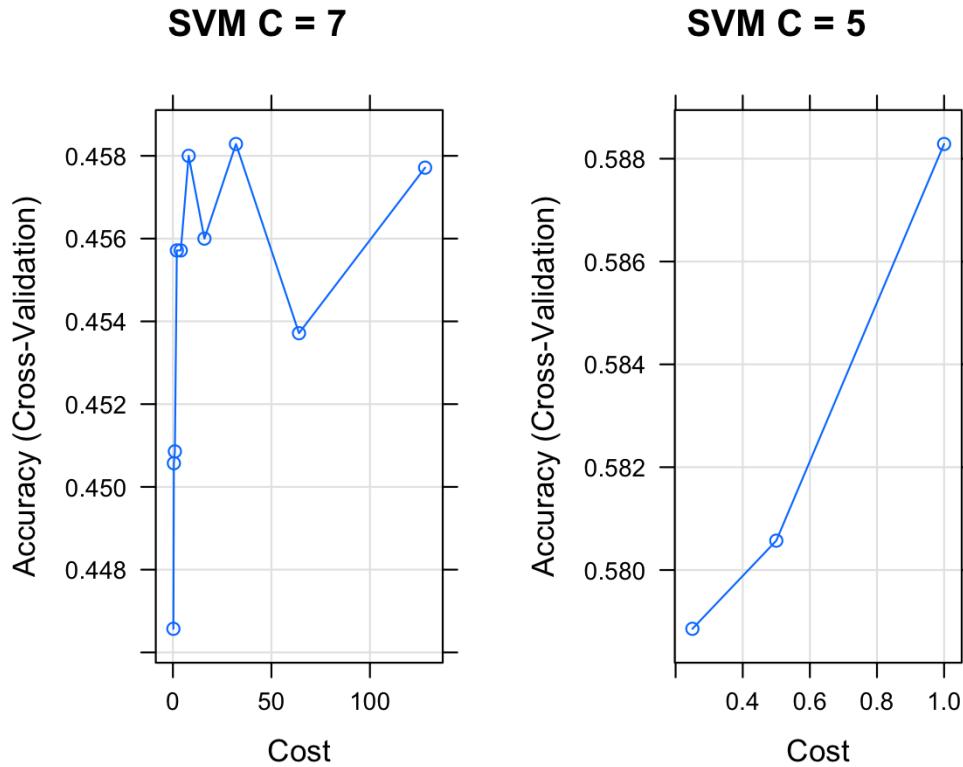


Figure 2.5: Comparison of accuracy results from different cost parameters for radial basis kernel function.

Polynomial Support Vector Machine

Figure 2.6 displays the results of different cost parameters on training accuracy for the polynomial basis function kernel. The final values used for the model predicting 7 classes were degree = 3, scale = 0.1, and cost = 1, which corresponded to a training accuracy of 0.44. The final model had a total of 2949 support vectors. The final model predicting only 5 classes produced an accuracy of 0.58 under degree = 3, scale = 0.1, and cost = 0.5. This final model had 2659 support vectors.

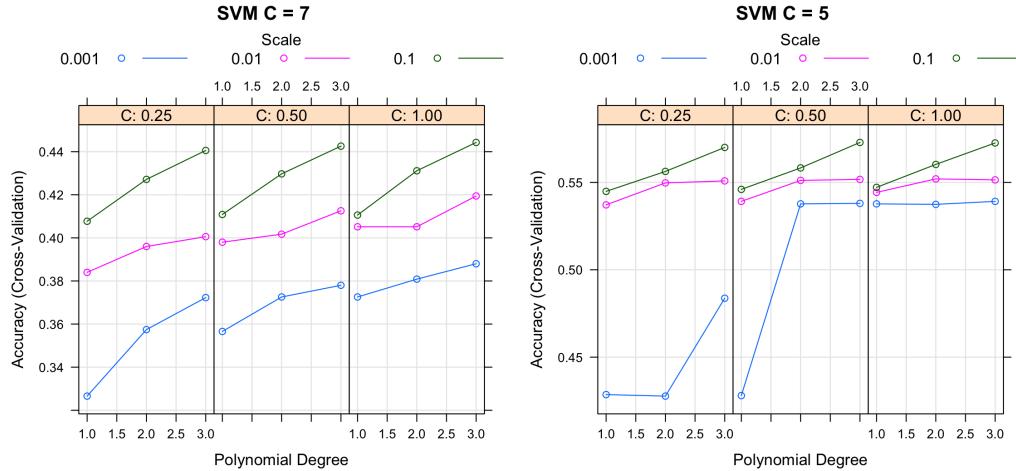


Figure 2.6: Comparison of accuracy results from different cost parameters with scale ranging from 0.001 to 0.1, C from 0.25 to 1, and degree from 1 to 3 for polynomial SVM.

Best Support Vector Machine

Figure 2.7 displays the results of the linear, radial basis, and polynomial basis support vector machine models. Based on the overall prediction accuracy, the radial basis is an appropriate kernel choice for the pixels data with the full 7 class prediction as well as 5 class prediction since it has a higher mean and median model accuracy than the other two kernel options.

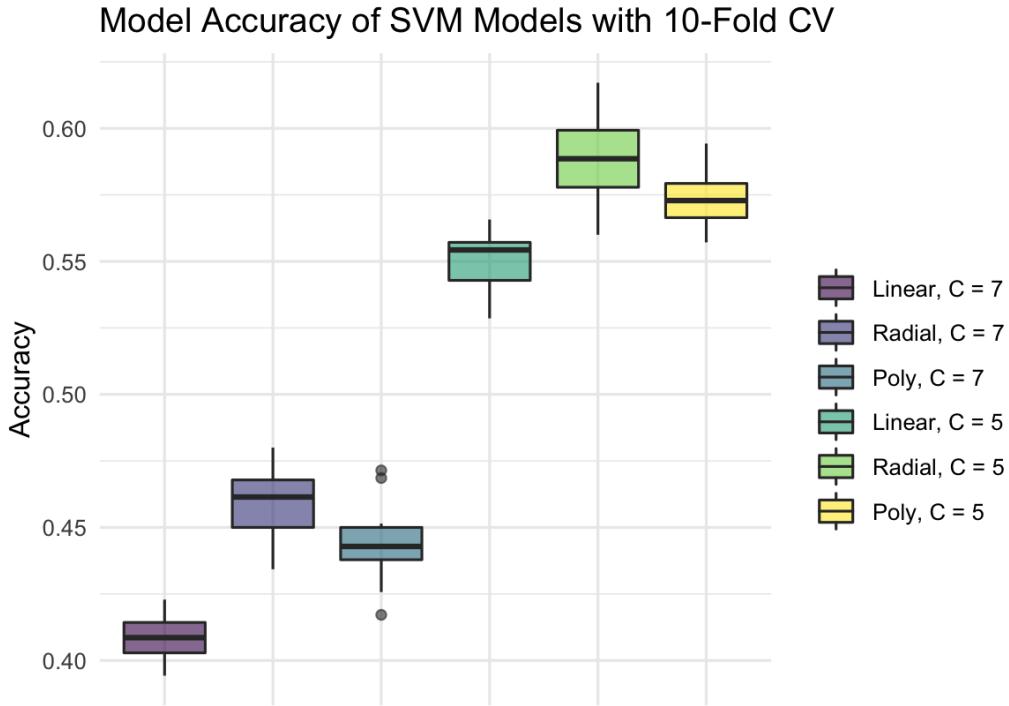


Figure 2.7: Comparison of results from support vector machine models with different kernel types: linear, radial, and polynomial.

As with the random forest model comparisons, a confusion matrix was constructed using the average counts over all ten cross-validation resamples to obtain the model results predicting Western redcedars. The support vector machine model that performed best overall was the radial basis kernel model, and it had the highest prediction accuracy of the Western redcedar class, 36%. The polynomial basis kernel had a 34% prediction accuracy and the linear kernel had 33% accuracy for Western redcedar. For Western redcedar prediction on the grouped classes, the radial model had 46% accuracy, and the polynomial model had 33% accuracy, and the linear model did not predict any Western redcedar pixels on average.

Table 2.2 compares the hyperparameters across all support vector machine models.

2.3 Testing Models

A subset of the data was withheld from training the models as test data to provide some measure of how the models perform in predicting results that the answer is unknown. Based on overall and redcedar accuracy, the appropriate models to consider with the testing data are the 7 class 8 predictor random forest, the 7 class radial support vector machine, the 5 class 8 predictor random forest, and the 5 class polynomial support vector machine. The results for the overall test accuracy and the prediction test accuracy for Western redcedars are shown in Tables 2.3 and 2.4. According to

Table 2.2: Accuracy of support vector machine models with different kernels. Best parameter is determined by ten-fold cross validation.

SVM Model Kernel	Accuracy	Kappa
Linear, C = 7	0.4080	0.3093
Radial, C = 7	0.4583	0.3680
Poly, C = 7	0.4443	0.3517
Linear, C = 5	0.5503	0.2817
Radial, C = 5	0.5883	0.3716
Poly, C = 5	0.5729	0.3295

Table 2.3: Overall accuracy of 8 predictor RF (Full), Radial SVM (Full), 8 predictor RF (Grouped), and Polynomial SVM (Grouped) models on test data

Type	RF Accuracy	SVM Accuracy
Full	0.5346	0.5318
Grouped	0.6497	0.6479

Table 2.4: Western redcedar accuracy of 8 predictor RF (Full), Radial SVM (Full), 8 predictor RF (Grouped), and Polynomial SVM (Grouped) models on test data

Type	RF Accuracy	SVM Accuracy
Full	0.3356	0.3056
Grouped	0.2407	0.0069

Table 2.5: Accuracy of radial SVM model with 7 classes and 8 predictor RF model with 5 classes for polygons in test data. A 'correct' prediction is considered to be a polygon with more than half of the pixels correctly classified.

Type	Result	n
Full SVM	Correct	199
Full SVM	Incorrect	474
Grouped RF	Correct	353
Grouped RF	Incorrect	320

the test data, the best performing models for further analysis are the radial support vector machine model for 7 class prediction and the 8 predictor random forest model for 5 class prediction.

The polygon prediction accuracy on the test dataset was also computed to see how pixels of the same tree were classified compared to individual pixel classification. A "correct" prediction was indicated if more than half the pixels in a polygon are correctly predicted. Table 2.5 displays the results over polygons in the test dataset.

Chapter 3

Results

3.1 Training Results

Predicting on the grouped data produced the highest overall accuracies. The best model on the grouped data was the random forest training model with an `mtry` value of 6 predictors at each split and the following predictors: red, green, blue, infrared, NDVI, $\frac{\text{red}}{\text{green}}$, $\frac{\text{red}}{\text{blue}}$. This model had a training predictive accuracy of 0.57 overall and 0.37 for Western redcedars specifically. On the test dataset, this model accurately predicted 24% of the pixels, while the accuracy for predicting more than half the pixels correctly in the same polygon was -%.

The model that performed best at predicting all 7 tree classes was the RF `P = 8` model with `mtry 2`. This model had a training accuracy of 0.44 overall, 0.33 for just Western redcedars, and testing accuracy of 0.53 overall and 0.34 for Western redcedar pixels. For polygon predictions, the RF `P = 8` model had an accuracy of -%. Between the RF `P = 7` model on the 5 classes data and the RF `P = 8` model on the 7 classes data, the RF `P = 7` model outperforms the RF `P = 8` model. These models are used to proceed with the final model over Portland.

3.2 Modelling Tree Species in Portland

After masking the raster images of the entire region of Portland, the pre-trained models are used to predict the tree species of individual pixels in the masked raster images. The `predict.raster` function from RStudio's `raster` package takes a model and applies that model to each cell of the raster, leaving a raster with tree classification predictions. This function is used to predict tree species over all five raster strips, and the raster strips are merged to display predictions for the entire city. Figure 3.1 displays the tree classification predictions given by the optimal random forest model over Portland and figure 3.2 displays the predictions from the support vector machine model.

RF Western Redcedar Predictions

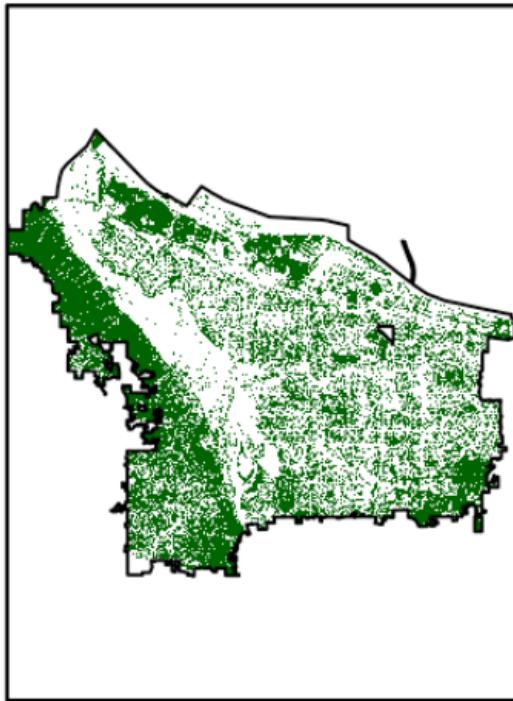


Figure 3.1: Model tree classification predictions of Western Redcedars over entire Portland region using the random forest model with 8 predictors on 7 classes.

RF Grouped Western Redcedar Predictions

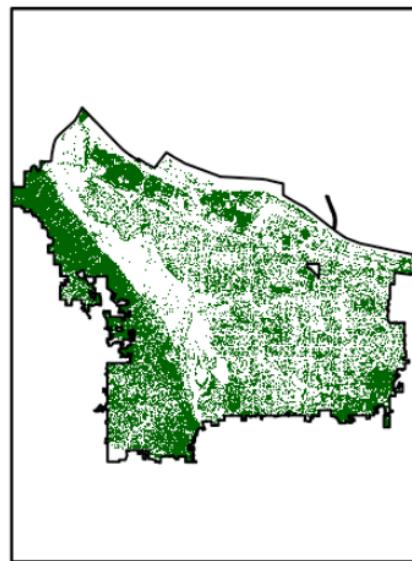


Figure 3.2: Model tree classification predictions of Western Redcedars over entire Portland region using RF model with 7 predictors on 5 classes.

In comparing the RF $C = 5$ model results to the RF $C = 7$ results, 83% of the pixel predictions were the same predictions under both models.

To assess the performance of the model over the Portland region, the results are compared to the remaining Western redcedars from `pdxTrees` both street and park data. The shapefiles for `pdxTrees` Western redcedars ground data are mapped over the classified pixel predictions, and the number of correctly and incorrectly classified pixels is recorded to be 52% accurate for RF $C = 7$ and 66% for RF $C = 7$. Figure 3.3 visualizes this comparison.

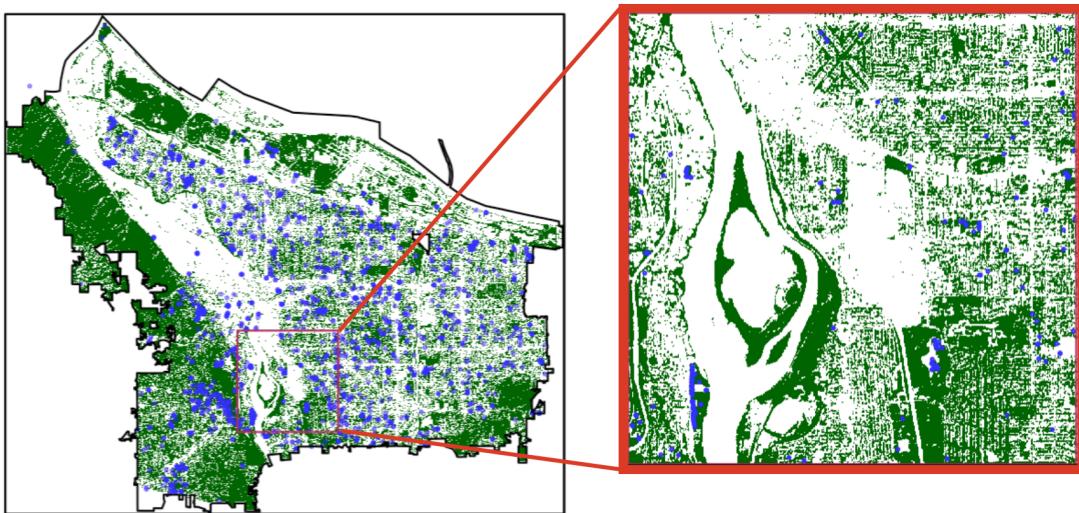
RF C = 7 with pdxTrees

Figure 3.3: Comparison of Western redcedar RF pixel predictions (in green) to Western redcedar trees from pdxTrees (in blue).

Chapter 4

Discussion

4.1 RF Grouped Data Performance Versus SVM Full Data Performance

The RF model predicting 5 classes outperforms the SVM model predicting all 7 classes in training and testing overall accuracy as well as in prediction of the whole polygons. It was observed that the models predicting less classes (Broadleaf, Douglas-Fir, Giant Sequoia, Grass, and Western Redcedar) outperformed models predicting on a larger set of tree classes (Bigleaf Maple, English Oak, Norway Maple, Douglas-Fir, Giant Sequoia, Grass, and Western Redcedar). This is likely a result of only having 4-band imagery plus the variables that are linear combinations of these 4 bands; 8 predictors training a model to predict 7 classes will not predict as well as 8 predictors training a model to predict 5 classes. In comparing the entire RF results on Portland to the `pdxTrees`, the test accuracy was 0.65.

4.2 Model Results Compared to Fricker Methods (Fricker et al., 2019)

Fricker's methods are conducted under ideal circumstances. The single image strip used in the study had a spatial resolution of 1 meter, they had access to both hyperspectral imagery and RGB imagery, and the ground data was carefully collected to consist of seven tree species (Fricker et al., 2019). Fricker's results show that models using hyperspectral images performed better than RGB image models because they provide additional information by including spectral band combinations not available from RGB images. For this research, the images only consist of 4 bands (red, green, blue, and infrared), are accessible online, and have a spatial resolution of 3 meters. Fricker's methods applying a convolution neural network using hyperspectral imagery produced cross-validation training accuracies of 87%. On just RGB imagery, this accuracy dropped to 64%, which is slightly less than the best performing model (RF P = 8, C = 5) from this research which had a cross-validation training accuracy of 65%.

4.3 Methods of Improvement/Further Work

As Fricker's article has shown, increasing the spectral resolution by having access to hyperspectral data adds more information to the models and improves performance. Hyperspectral imagery collects information from more regions of the spectrum than RGB imagery, which allows it to capture finer details. Tree and vegetation pixels reflect a certain amount of light, especially regions of dense forests, so having more bands can allow the models to detect subtle differences in vegetation pixels with non-vegetation pixels. By increasing the number of variables in the model, the risk of multicollinearity that comes from artificially creating predictors with linear combinations of the RGB bands also gets removed. The benefits of high spatial and spectral resolution imagery also extends to improving the masks applied to the images, which were successful, for example, in removing building structures and dead fields of grass, but incorrectly recognized mossy lakes as trees. Carefully masking the images improves the ability of the model to locate tree pixels and classify them.

In terms of data, another improvement to this research would be to increase the spatial resolution to 1 meter or better if possible. This refines the process of creating polygons by enhancing the visibility of the tree canopy outlines and reduces noise introduced by having shadows, dirt, building structures, and other non-tree pixels within the polygons. Another method to improve the results is by improving the drawn polygons themselves to only include pixels that represent the tree canopy. Drawing the polygons for each tree species by hand is likely to introduce error to the analysis because the first few polygons are not likely to be drawn the same way that the last few polygons are drawn. This issue is addressed by taking care to only include pure tree pixels (no shadows, dirt, or other structures) and ensuring that each polygon consists of at least 10 pixels.

An ideal training data set for the models would consist of data from imagery with high spatial and spectral resolution and also polygons with pure tree canopy pixels. If multiple raster strips are necessary for the area of interest, the polygons need to be evenly distributed across all raster strips and ideally the tree species per polygon as well. The trees themselves should have large canopies that hopefully do not overlap with other tree canopies of different species.

With these improvements, the research will allow researchers to have a better understanding of the locations of redcedars in Portland. From there, this information provides a way to track future changes in redcedar trees's health, and inform decisions about tree management in the city. In a broader context, applying this method to Portland imagery data with multiple strips provides information about the applicability of modelling tree locations in different parts of the world and with data from different satellites. Ultimately, tracking the changes of the ecosystem allows people to take steps to preserve dying species.

4.4 Conclusion

Reports of Western redcedars in the Pacific Northwest over the past 10 years have sparked an interest in investigating the decline of the species. This research combined RGB imagery and ground-level data available for the city of Portland to try and predict the locations of Western redcedars and progress towards understanding the decline of the species and eventually expand the methods to larger regions like the Pacific Northwest. Ground-level data was used to locate individual tree canopies and label their species on the satellite images, then the light intensities were extracted for each tree canopy pixel into a training data set. Random forest and support vector machine models were trained to classify 7 or 5 tree species. The best performing model was the random forest model with 7 predictors on 5 tree species with an F-score of 0.5 for Western redcedars. These results were not as high as other results from data with better conditions (higher spatial and spectral resolutions), however, they provide a better way of classifying a tree pixel as Western redcedar than a random guess of one of the 7 species (probability of $1/7 \approx 0.14$).

The results of this research are replicable, since it involves data from open sources and the coding is completed in RStudio, and not too computationally intensive to require more than a day to run all of the code. Another key takeaway is that this research applies the methods of using remote sensing imagery to classify tree species in Portland specifically and provides a guideline for general areas to focus on locating Western redcedars. The research is another application of remote sensing for tree species classification, which has the potential to predict more tree species in larger regions with more research.

Appendix A

The First Appendix

This first appendix includes all of the R chunks of code that were hidden throughout the document (using the `include = FALSE` chunk tag) to help with readability and/or setup.

In the main Rmd file

In Chapter ??:

Appendix B

The Second Appendix, for Fun

References

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