Neotoma paper

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# neotoma: A Programmatic Interface to the Neotoma Paleoecological Database

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## Abstract:

Paleoecological data are integral to ecological analyses. First, they provide an opportunity to study ecological and evolutionary interactions between communities and abiotic environments over long time scales. Second, they allow us to study processes that occur infrequently, such as megadroughts, hurricanes, and rapid climate change. Third, the past allows us to study ecological processes in the absence of widespread anthropogenic influence.

The R package neotoma, described here, obtains and manipulates data from the Neotoma Paleoecological Database (Neotoma Database: <http://www.neotomadb.org>). The Neotoma Database is a public-domain searchable repository for multiproxy paleoecological records spanning the past 5 million years. The Neotoma Database provides the cyberinfrastructure to study spatiotemporal dynamics of species and communities from the Pliocene to the present; neotoma provides a user interface to enable these studies. The neotoma searches the Neotoma Database using search keys that can include location, taxon name, or dataset type (e.g., pollen, vertebrate fauna, ostracode) using the Database's Application Programming Interface (API). The package returns a set of nested metadata associated with the site, including the full assemblage record, geochronological data to enable rebuilding of age models, metadata for the dataset (e.g. age range of samples, date of accession into Neotoma, principal investigator), and site metadata (e.g. location, site name and description). neotoma also provides tools to allow cross-site analysis, including the ability to standardize taxonomies using built-in taxonomies derived from the published literature or user-provided taxonomies.

To assist with the use of the neotoma package we provide examples of key functions based on the published literature, for both plant and mammal taxa.

Keywords: R software, neotoma, paleoecology, database, pollen, mammal

## Introduction

Paleoecological data are fundamental to understanding the patterns and drivers of biogeographical, climatic, and evolutionary change, ranging from the recent past to the dawn of life. Although individual site-level studies have provided fundamental insights into past ecological dynamics, the true power of paleoecological data emerges from networks of paleoecological data assembled to study broad-scale ecological and evolutionary phenomena, e.g. the responses of speciation rates to the five major extinction events in geological history (Peters & Foote 2001; Raup & Sepkoski 1984; Sepkoski 1997) and the rapid and individualistic responses of species to the climate changes accompanying recent glacial-interglacial cycles (Davis 1981; Schroeder et al. 1996; Huntley & Webb 1988; Tzedakis 1994; Williams et al. 2004). Paleoecoinformatics (Brewer et al. 2012; Uhen et al. 2013) is dedicated to providing tools to researchers across disciplines to access and use large paleoecological datasets spanning thousands of years. These datasets may be used to provide better insight into regional vegetation change (Blois et al. 2013; Blarquez, Carcaillet, et al. 2014), patterns of biomass burning (Marlon et al. 2013), or changing rates of geophysical processes through time (Goring et al. 2012). The increasing interest in uniting ecological and paleoecological data to in order to better understand responses to a rapidly changing world (Fritz et al. 2013; Behrensmeyer & Miller 2012; Dietl & Flessa 2011) will require more robust tools to access and synthesize data from the modern and paleo time domains.

The Neotoma Paleoecological Database represents a consortium of paleoecological databases, with distributed scientific governance and expertise, but sharing a common database infrastructure. Constituent databases include, among others, the large European, Latin American, and North American Pollen Databases; the North American Plant Macrofossil Database; FAUNMAP (Pliocene to Quaternary mammal fossils in the United States and Canada); the North Dakota State University Fossil Insect Database; the North American Non-Marine Ostrocode Database; and the Diatom Paleolimnology Data Cooperative. Neotoma is the outgrowth of a longstanding collaboration between the European Pollen Database and the North American Pollen Database (Grimm et al. 2013) and the desire to integrate these data with faunal and other paleo data. The database framework was generalized from the pollen databases (which had identical structures) and the FAUNMAP database to accomodate both macro- and microfossil data as well as other kinds of data such as geochemical, isotopic, and loss-on-ignition. Work is underway to include other taxonomic groups and depositional contexts (e.g. testate amoeba records, packrat midden data), thus further expanding the data that can be accomodated by Neotoma. Crucially, Neotoma is a vetted database. Through the use of data stewards - domain experts distributed among constituent bases who can check for inaccuracies, upload and manage data records - Neotoma can support high qualitycontrol assurance for each of the constituent data types, and receive feedback from research communities involved with each specific data type (Grimm et al. 2013).

The Neotoma Database has also developed Application Programming Interfaces that allow users to query the database via web services, which return data using properly formed URL requests. For example, the URL: <http://api.neotomadb.org/v1/apps/geochronologies/?datasetid=8> will return all geochronological data for the record associated with the dataset ID 8.

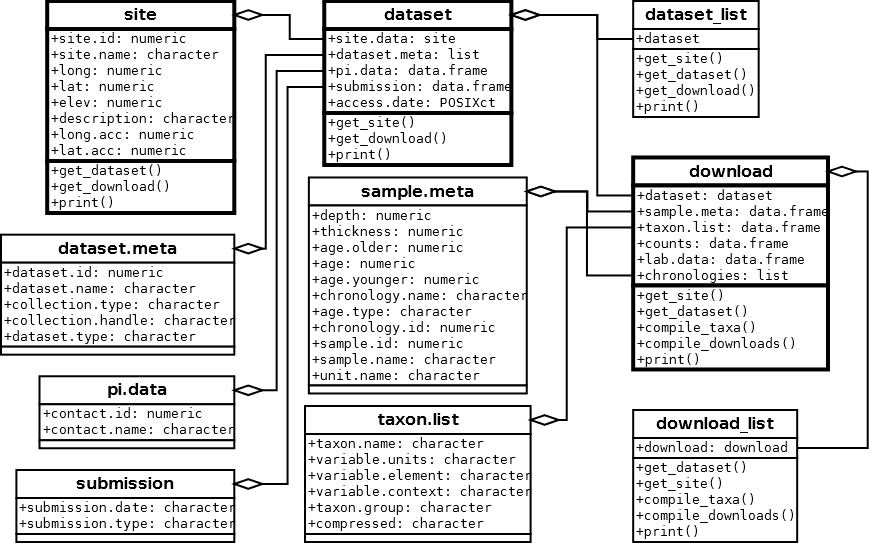
The analysis of paleoecological data is commonly performed using the statistical software R (R Core Team 2014), and several paleoecological packages in R packages are designed specifically for paleoecological dataexist, for analysis including analogue (Simpson & Oksanen 2014; Simpson 2007) and rioja (Juggins 2013) for paleoenvironmental reconstruction, Bchron for radiocarbon dating and age-depth modeling (Parnell 2014) and paleofire to access and analyse charcoal data (Blarquez, J. R. Marlon, et al. 2014). Given the rapid proliferation and availability of these analytical tools in R, the rate-limiting step has become the difficulty of obtaining and importing data into R. This bottleneck has meant reliance on datasets such as those from the NOAA Paleoclimate Repository or the North American Modern Pollen Database, and on more *ad hoc* methods such as the distribution of individual datasets from author to analyst.

With an increasing push to provide ecological publications that include numerically reproducible results (Goring et al. 2013; Goring et al. 2012; Wolkovich et al. 2012; Reichman et al. 2011) it is important to provide tools that allow analysts to directly access dynamic datasets, and to provide tools to support reproducible workflows. The rOpenSci project is dedicated to developing tools using R to facilitate a culture shift toward reproducible science in the ecology comunity. As part of this effort, it has provided a number of tools that can directly interact with application programmatic interfaces (APIs) to access data from a number of databases including rfishbase for FishBase (Boettiger et al. 2012), and taxize for the Encyclopedia of Life, iPlant/Taxosaurus and others (Chamberlain & Szöcs 2013) among others.

The neotoma package addresses concerns regarding data access and workflow reproducibility by providing users with tools that allow paleoecologists to query, download, organize, and summarize data from the Neotoma database using R. Here we describe the neotoma package, then we present use cases for the neotoma package, using examples drawn from the ecological literature, with the general objective of illustrating how neotoma provides tools to perform paleoecological research in an open and reproducible manner.

## The neotoma package

neotoma R package is an interface between the Neotoma Paleoecological Database (<http://neotomadb.org>) and statistical tools in R. neotoma uses an API to send data requests to the Neotoma Database, and then forms data objects that can interact with existing packages such as analogue (Simpson & Oksanen 2014) and rioja (Juggins 2013), which are used for environmental reconstruction, manipulation, and presentation of paleoecological data. The neotoma package also includes tools to standardize pollen taxon names across sample sites using a set of commonly accepted pollen taxonomies for North America, or user defined taxonomies.

 **Figure 1**. *Major classes in neotoma, their relations to one another and the associated functions. The classes described below have a heavier outline than their associated variables.*

Data in the neotoma package is represented in three main classes (Figure 1): "site"s, "dataset"s (grouped into "dataset\_list"s), and "download"s (grouped into "download\_list"s). A "site" is the most basic form of spatial information representing the spatial locations of datasets along with site names, descriptions and a unique site.id. "site"s are "data.frame"s with columns siteid, sitename, lat, long, elev, description, long\_acc, and lat\_acc. These column headings are generally self explanatory; long\_acc and lat\_acc are used to indicate the width of the bounding box for a sample site (with a midpoint of long and lat). In the Neotoma Database, examples of sites include a lake from which one or more cores are collected, a cave from which one or more faunal assemblages are collected, an archaeological dig with one or more excavation pits, and so forth. Each unique site returned by the get\_site() search is placed in a row, providing enough descriptive data to plot locations and understand the spatial context of a site.

Although get\_site() is useful for first-pass surveys of data availability, analysts more commonly will want to to search for and retrieve datasets stored in the Neotoma Database. The neotoma package allows you to use almost all of the same search terms in get\_dataset() as in get\_site(), and returns a more complete description of the datasets available, however at this time get\_site() is the only method by which you can search for site names.

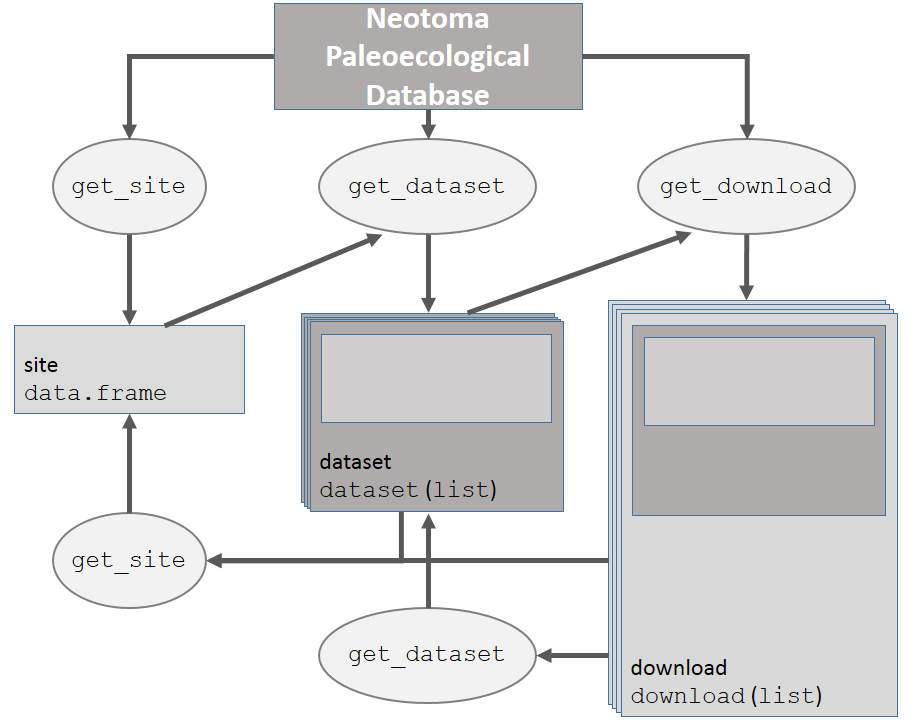
"dataset"s associated with individual sites can be obtained using get\_dataset(). Each search using get\_dataset() returns a "dataset\_list", a list of "dataset"s, equal to the number of datasets returned by the function. A dataset in the Neotoma Database is a set of samples of the same type from a single collection unit within a site. Examples of datasets in Neotoma include 1) all the pollen counts from a single core from a lake, 2) all the geochronological measurments (e..g radiocarbon dates) from a sediment core, 3) all the faunal data from an excavation in a cave, 4) all the plant macrofossil data from a packrat midden . In neotoma, a "dataset" includes the "site" for each "dataset" (as "site.data", Figure 1), along with the "dataset" specific metadata ("dataset.meta" in Figure 1). The dataset also includes the principal investigator, submission date to Neotoma, and the date that the information was accessed from the Neotoma API using the R package. "dataset\_list"s and "dataset"s can be used to access the full "download" using the get\_download() method.

get\_download() returns an object of class "download\_list" containing objects of class "download" (Figure 1). get\_download() will also return a confirmation message for each individual API call as the function proceeds that can be turned off using the argument verbose = FALSE. Each "download" contains the associated "dataset" (Figure 1). The "sample.meta" component is where the core depth and age information is stored. The actual chronologies are stored in "chronologies". If a core has a single age model then "chronologies" has a length of one. Some cores have multiple chronologies and these are added to the list. The default chronology is always represented in "sample.meta" and is always the first chronology.

To build a new chronology with the same chronological controls as an existing chronology, but with a different algorithim, use get\_chroncontrol() to return the chronology controls and the "chronology.id" in either "sample.meta" or any one of the "chronologies" objects. While the chronological controls used to build a chronology may vary across chronologies for a single site, the default model contains the "best" chronological control data, as determined at the time the chronologies for the collection unit were last reviewed. It is important to note, however, that the "best" chronologies for most collection units in the database were based on "classical" age models (Blaauw 2010) that do not include estimates of uncertainty. Moroever, these default age models that are in calibrated radiocarbon years utilize radiocarbon dates that are calibrated a priori. Bayesian age modeling programs, such as Bacon (Blaauw & Christen 2011), which provides estimates of uncertainty, as well as the classical age modeling program clam (Blaauw 2010), which also returns estimates of uncertainty, utilize uncalibrated radiocarbon dates as input. Thus, the calibrated ages of many existing age models in Neotoma will not be appropriate for these programs, and the age controls may have to be obtained from the geochronology table using get\_geochron().

The age controls of existing default radiocarbon-year chronologies may sometimes be appropriate for programs such as Bacon and clam; however, many of these chronologies rejected radiocarbon dates *a priori*, which could be included when using Bacon. The neotoma package has a function to interface directly with Bacon or clam, called write\_agefile(), which will output a correctly formatted age file for either of these applications using a "download" object.

The "taxon.list" lists taxa found in the dataset samples ("counts") and any laboratory data ("lab.data"), along with the units of measurement and taxonomic grouping (Figure 1). The counts are the actual count, presence or percentage data recorded for the dataset. The lab.data component contains information about any spike used to determine concentrations, sample quantities and, in some cases, charcoal counts.

 **Figure 2**. *How the main data objects relate to one another in the neotoma package, and the helper functions used to move from one data type to another.*

Each of these objects, "site", "dataset" and "download" can be obtained using direct calls to the API, or using functions defined in the neotoma package (Figure 2).

## Examples

Here we present several examples that both introduce users to the neotoma package, and highlight how neotoma can be used in a paleoecological worklfow. We beging with a simple example in which we compare change in *Alnus* pollen between two sites during the late Quaternary, followed by two more involved examples where we look at *Pinus* migration and changes in late-Quaterary mammal distributions in the United States.

### A simple example

A researcher is interested in finding the pollen record for Marion Lake, in British Columbia (Mathewes 1973) and comparing changes in *Alnus* pollen to the Louise Pond record (Pellatt & Mathewes 1997) from Haida G'Waii, further north. We search for specific sites by name using get\_site(), making use of the wildcard "%" to catch sites whose site names begin with Marion Lake or Louise Pond:

library("neotoma")  
library("analogue")  
  
marion <- get\_site(sitename = "Marion Lake%")

The API call was successful, you have returned 1 records.

louise <- get\_site(sitename = "Louise Pond%")

The API call was successful, you have returned 1 records.

louise

siteid long lat elev  
 Louise Pond 1618 -131.8 53.42 650  
 description  
 Louise Pond Glacial scour lake. Physiography: Queen Charlotte Ranges, Louise Island. Surrounding vegetation: Tsuga mertensiana, Pinus contorta.  
 long.acc lat.acc  
 Louise Pond 0 0

In each case get\_site() returns a single "site" (Figures 1 & 2). Here we queried the Neotoma database for site based on sitename, but alternately we could have queried for sites within a geographical bounding box, or by geopolitical region.

To get "dataset"s for these records we can rbind()ing the two records, and use get\_dataset() directly (Figure 2):

western.sites <- rbind(marion, louise)  
western.data <- get\_dataset(western.sites)

"western.data" is a "dataset\_list", containing two "dataset"s (Figure 1). The "dataset" for a site will be nested within a "dataset\_list", even if only a single site is returned, so that methods can be consistent across classes and functions. This means that a single "dataset" must be retrieved as *e.g.*, western.data[[1]] (this is also the case for "download" and "download\_list" objects). The use of "dataset" and "dataset\_list" classes allow us to easily move between get\_dataset(), get\_site() and get\_download(). We can see the special print() method for both datasets and dataset\_lists before we download the full records and print them:

western.data

A dataset\_list containing 2 objects:  
 Accessed from 2014-10-05 15:09h to 2014-10-05 15:09h.   
 Datasets:  
 dataset.id site.name long lat type  
 1705 Marion Lake (CA:British Columbia) -122.5 49.31 pollen  
 1670 Louise Pond -131.8 53.42 pollen

western.data[[1]]

A dataset for Marion Lake (CA:British Columbia)  
 Accessed 2014-10-05 15:09h.   
 dataset.id site.name long lat type  
 1705 Marion Lake (CA:British Columbia) -122.5 49.31 pollen

western.dl <- get\_download(western.data)

API call was successful. Returned record for Marion Lake(CA:British Columbia)  
 API call was successful. Returned record for Louise Pond

Warning:   
 Modifiers are absent from the lab objects Lycopodium tablets, Lycopodium spike, Sample quantity.   
 get\_download will use uniqueidentifiers to resolve the problem.

western.dl

A download\_list containing 2 objects:  
 Accessed from 2014-10-05 22:16h to 2014-10-05 22:16h.   
 Datasets:  
 dataset.id site.name long lat age.younger  
 1705 Marion Lake (CA:British Columbia) -122.5 49.31 58  
 1670 Louise Pond -131.8 53.42 62  
 age.older type  
 13051 pollen  
 10065 pollen

western.dl[[1]]

A download object for Marion Lake (CA:British Columbia)  
 Accessed 2014-10-05 22:16h.   
 dataset.id site.name long lat age.young  
 1705 Marion Lake (CA:British Columbia) -122.5 49.31 58  
 age.old type  
 13051 pollen

Pollen taxonomy can vary substantially across cores depending on the level taxonomic resolution used by a pollen analyst, or because of changes to taxonomies over time. One analyst might discriminate subgenera of *Pinus*, another might simply identify *Pinus* to the genus level. Gramineae - a common pollen type in earlier pollen records - has now been renamed Poaceae. This variable and shifting taxonomy is a first-order challenge for analysts seeking to analyze the dynamics of taxa across multiple groups. neotoma provides several options for standardized taxonomic list, corresponding to three published taxonomies for the United States and Canada (Gavin et al. 2003; Whitmore et al. 2005; Williams & Shuman 2008). This function can be helpful, but it should also be used with care. The aggregation table is accessible using the command data(pollen.equiv) and the function to compile the data is called compile\_taxa(). It can accomodate either the internal translation table provided with the package, or a user-defined table with the same structure as pollen.equiv.

We are interested in comparing the relative pollen abundances of a *Alnus* between two sites. We compile the pollen data using the 'P25' taxonomy from Gavin et al. (2003). The first record downloaded is Marion Lake. We can see the "download" for Marion Lake the taxon.table has 5 columns:

head(western.dl[[1]]$taxon.list)

|  |  |  |
| --- | --- | --- |
| taxon.name | variable.units | variable.element |
| Tsuga heterophylla | NISP | pollen |
| Poaceae | NISP | pollen |
| Unknown (monolete) | NISP | spore |
| Tsuga mertensiana | NISP | pollen |
| Rosaceae | NISP | pollen |
| Salix | NISP | pollen |

Table continues below

|  |  |
| --- | --- |
| variable.context | taxon.group |
|  | Vascular plants |
|  | Vascular plants |
|  | Unidentified palynomorphs |
|  | Vascular plants |
|  | Vascular plants |
|  | Vascular plants |

Once we apply compile\_taxa() to the dataset using the 'P25' compiler:

western.comp <- compile\_taxa(western.dl, list.name = "P25")  
names(western.comp) <- c("marion", "louise")

The taxon.table for Marion Lake now has an extra column (note that several columns were removed to improve readability).

head(western.comp[[1]]$taxon.list[, c(1, 5, 6)])

|  |  |  |  |
| --- | --- | --- | --- |
|  | taxon.name | taxon.group | compressed |
| **2** | Tsuga heterophylla | Vascular plants | Tsuga |
| **29** | Poaceae | Vascular plants | Poaceae |
| **3** | Unknown (monolete) | Unidentified palynomorphs | Other |
| **4** | Tsuga mertensiana | Vascular plants | Tsuga |
| **5** | Rosaceae | Vascular plants | Other |
| **6** | Salix | Vascular plants | Salix |

compile\_taxa() returns a "download\_list" or "download", for which taxon.list gains a column named compressed to link the original taxonomy to the revised taxonomy. This linkage is an important reference for researchers who choose to use this package for large-scale analysis, but who might need to later check the aggregated taxonomic groups against the original data. In this example we see that all the spore types have been lumped into a single *Other*.

The sample data ("counts") contained in each "download" in the "download\_list" wester.dl are converted into percentages to standardize between cores using tran() from the analogue package (Simpson 2007) so we can see what happens with *Alnus* on the west coast of North America during the Holocene:

marion.alnus <- tran(x = western.comp$marion$counts, method = "percent")[, "Alnus"]  
louise.alnus <- tran(x = western.comp$louise$counts, method = "percent")[, "Alnus"]  
  
alnus.df <- data.frame(alnus = c(marion.alnus, louise.alnus), ages = c(western.comp$marion$sample.meta$age,   
 western.comp$louise$sample.meta$age), site = c(rep("Marion", length(marion.alnus)),   
 rep("Louise", length(louise.alnus))))  
  
plot(alnus ~ ages, data = alnus.df, col = alnus.df$site, pch = 19, xlab = "Years Before Present",   
 ylab = "Percent Alnus")

plot of chunk alnus-data-plot **Figure 3**. *Plots of* Alnus *pollen percentages at two sites, one in the lower mainland of British Columbia (Marion Lake) and the other on Haida G'waii (Louise Pond). Axis labels are presented as if the code was run directly, but represent calibrated radiocarbon years before present on the x-axis and* Alnus *pollen percentages on the y-axis.*

Marion Lake (red) maintains much higher proportions of *Alnus* throughout its history, and has a rapid increase in *Alnus* pollen during the historical period. This rapid shift in the last 200 years is likely as a result of rapid colonization by pioneer *Alnus rubra* following forest clearance and fire in the lower mainland of British Columbia (Mathewes 1973).

It is also possible to plot the pollen stratigraphy at any one site, again, using the analogue package for R (Simpson 2007). Here we plot Marion Lake:

core.pct <- data.frame(tran(western.comp[[1]]$counts, method = "percent"))  
  
core.pct$age <- western.comp[[1]]$sample.meta$age  
  
# Eliminate taxa with no samples greater than 4%.  
core.pct <- chooseTaxa(core.pct, max.abun = 4)  
  
Stratiplot(age ~ ., core.pct, sort = "wa", type = "poly")

plot of chunk marion-plot **Figure 4***.* Stratigraphic plot for Marion Lake. Age is plotted on the y-axis in calibrated radiocarbon years before present. The analogue package provides extensive opportunity to customize the stratigraphic plot beyond this simple example.\*

### Pinus migration following the last Glacial Maximum

Macdonald and Cwynar (1991) used pollen percentage data for *Pinus* to map the northward migration of lodgepole pine (*Pinus contorta* var *latifolia*) following the retreat of the Laurentide Ice Sheet and the accompanying rise of temperatures in the late-Glacial and early Holocene. In their study a cutoff of 15% *Pinus* pollen was defined as the indicator of *Pinus* presence. Strong and Hills (2013) have remapped the migration front using a lower pollen proportion (5%) and more sites. Here, the analysis is partially replicated. Note that additional R packages must be installed and loaded for the following examples.

A spatial bounding box delimiting sites is defined for the search. Strong and Hills (2013) use a region approximately bounded by 54oN and 65oN, and from 110oW to 130oW. The function get\_site() can return all sites within a slightly expanded bounding box:

# install.packages('ggmap', 'ggplot2', 'reshape2', 'Bchron', 'gridExtra')  
library("ggmap")  
library("ggplot2")  
library("reshape2")  
library("Bchron")  
library("gridExtra")  
  
all.sites <- get\_site(loc = c(-140, 45, -110, 65))

The API call was successful, you have returned 444 records.

The code above returned 444 sites. To narrow down the search we will use get\_dataset() to search for all Pinus taxa within the same bounding box as above. get\_dataset() can also limit the type of dataset, either by looking for specific taxa, or by describing the dataset type (e.g., datasettype = 'pollen' or datasettype = 'mammal'). The % wildcard indicates that any characters may follow a string starting with "Pinus":

all.datasets <- get\_dataset(loc = c(-140, 45, -110, 65), datasettype = "pollen",   
 taxonname = "Pinus%")

The API returns 69 datasets. Many dropped sites were pollen surface samples, or sites with datasets for other taxonomic groups. The distribution of the 69 fossil pollen sites can be plotted over our original 444 sites. We use ggplot2 (Wickham 2009) to make the figures:

map <- map\_data("world")  
ggplot(data = data.frame(map), aes(long, lat)) + geom\_polygon(aes(group = group),   
 color = "steelblue", alpha = 0.2) + geom\_point(data = all.sites, aes(x = long,   
 y = lat)) + geom\_point(data = get\_site(all.datasets), aes(x = long, y = lat),   
 color = 2) + xlab("Longitude West") + ylab("Latitude North") + coord\_map(projection = "albers",   
 lat0 = 40, lat1 = 65, xlim = c(-140, -110), ylim = c(45, 70))

plot of chunk map-pinus-example **Figure 5** *Mapped sites with pollen cores in the interior of British Columbia and the Yukon Territory of Canada (red), including other Neotoma sites without stratigraphic pollen data (black).*

The map (Figure 5) shows a number of sites in the interior of British Columbia that have no fossil pollen. There are also other sites not shown here that may have relevant data but have yet to be entered into the database. This highlights a common challenge in paleoecoinformatics - the import of individual records into data repositories takes some time, and is an on-going process that is aided by the collective contributions of the original analysts, data stewards, and large-scale research initiatives (e.g. PAGES 2K, PalEON). Fortunately, new software tools are speeding up the process of uploading and vetting data. The Tilia software (<http://www.neotomadb.org/data/category/tilia>) has been updated to allow direct upload to the Neotoma Database and includes a large number of automated data quality checks and standardized look-up tables for variable names. Because neotoma directly links to the Neotoma Database via APIs, analyses using neotoma can be updated continuously as new sites are added.

To obtain the data for the 69 sites we use get\_download():

all.downloads <- get\_download(all.datasets, verbose = FALSE)

Only the percentage of *Pinus* is of interest, so we can again compile the taxa across the "download\_list" using the 'P25' taxonomy (Gavin et al. 2003):

compiled.cores <- compile\_taxa(all.downloads, "P25")

We want to determine which sample has the first local *Pinus* presence in each core using a cutoff of 5% (Strong & Hills 2013). We can find which rows in the *Pinus* column in each "download"'s "count" data.frame have presence over 5% and then find the highest row number since the samples in a dataset are ordered stratigraphically, with the youngest sample in the top row and the oldest sample in the bottom row. This is a more complicated example:

top.pinus <- function(x) {  
   
 x.pct <- tran(x$counts, method = "proportion")  
   
 # Cores must span at least the last 5000 years (and have no missing dates):  
 old.enough <- max(x$sample.meta$age) > 5000 & !all(is.na(x$sample.meta$age))  
   
 # Find the highest row index associated with Pinus presence over 5%  
 oldest.row <- ifelse(any(x.pct[, "Pinus"] > 0.05 & old.enough), max(which(x.pct[,   
 "Pinus"] > 0.05)), 0)  
   
 # return a data.frame with site name & location, and the age and date type  
 # (since some records have ages in radiocarbon years) for the oldest Pinus.  
   
 if (oldest.row > 0) {  
 return(data.frame(site = x$dataset$site.data$site.name, lat = x$dataset$site.data$lat,   
 long = x$dataset$site.data$long, age = x$sample.meta$age[oldest.row],   
 date = x$sample.meta$age.type[oldest.row]))  
 } else {  
 NULL  
 }  
}  
  
# Apply the function 'top.pinus' to each core using lapply and rbind:  
summary.pinus <- do.call(rbind.data.frame, lapply(compiled.cores, top.pinus))

We need to calibrate dates that are recorded in radiocarbon years using the Bchron package (Parnell 2014). In most cases the original uncertainty for individual ages is not recorded in "sample.meta" so we assume a 100 year standard deviation, however, direct recalibration of radiocarbon dates from interpolated age models is not the best approach, but we use it as an example here:

radio.years <- summary.pinus$date %in% "Radiocarbon years BP"  
  
# BChronCalibrate is in the BChron package:  
calibrated <- BchronCalibrate(summary.pinus$age[radio.years], ageSds = rep(100,   
 sum(radio.years, na.rm = TRUE)), calCurves = rep("intcal13", sum(radio.years,   
 na.rm = TRUE)))  
  
# we want the weighted means from 'calibrated'  
wmean.date <- function(x) sum(x$ageGrid \* x$densities/sum(x$densities))  
  
summary.pinus$age[radio.years] <- sapply(calibrated, wmean.date)  
summary.pinus <- na.omit(summary.pinus)  
  
summary.pinus <- summary.pinus[!((summary.pinus$age < 2000) & (summary.pinus$long <   
 -130)), ]  
  
# A loess curve is straightforward, but not the best model:  
regress <- ggplot(summary.pinus, aes(x = lat, y = age)) + geom\_point(aes(color = age),   
 size = 2) + scale\_y\_reverse(expand = c(0, 100)) + xlab("Latitude North") +   
 ylab("Years Before Present") + geom\_smooth(n = 40, method = "loess") + geom\_rect(aes(xmin = 59,   
 xmax = 60, ymin = 7000, ymax = 10000), color = 2, fill = "blue", alpha = 0.01)  
  
mapped <- ggplot(data = data.frame(map), aes(long, lat)) + geom\_polygon(aes(group = group),   
 color = "steelblue", alpha = 0.2) + geom\_point(data = summary.pinus, aes(x = long,   
 y = lat, colour = age), size = 3) + coord\_map(projection = "albers", lat0 = 40,   
 lat1 = 65, xlim = c(-140, -110), ylim = c(40, 70)) + theme(legend.position = "none")  
  
grid.arrange(mapped, regress, nrow = 1)

plot of chunk pinus-recal-plot **Figure 6**. *Mapped ages of first* Pinus *establishment in the interior of British Columbia and the Yukon Territory based on a 5% pollen cut-off. The age of first appearance is also plotted and smoothed with a loess curve.*

The results show a clear pattern of northward expansion for *Pinus* in northwestern North America. These results agree broadly with the findings of Strong and Hills (2013) who suggest that *Pinus* reached a northern extent between 59oN and 60oN at approximately 10 - 7 ka cal BP as a result of geographic barriers before continuing northward after 7 ka cal BP.

### Mammal Distributions in the Pleistocene

Graham et al. (1996) built and applied the FAUNMAP dataset (<http://www.ucmp.berkeley.edu/faunmap/>) of fossil assemblages to elucidate patterns of change in mammal distributions through the Pleistocene to the present. The paper uses various multivariate analyses to show, in part, that mammal species have responded in a Gleasonian manner to climate change since the late-Pleistocene. Graham et al. (1996) show some species migrating northward in response to warming climates, others staying relatively stable, and some moving southward. FAUNMAP has been incorporated into Neotoma (and expanded with new records), and this example performs some simple analyses that show how different species responded to the changing climate of the last deglaciation.

First, all vertebrate fauna datasets are obtained from Neotoma:

# Bounding box is effectively the continental USA, excluding Alaska.  
mam.set <- get\_dataset(datasettype = "vertebrate fauna", loc = c(-125, 24, -66,   
 49.5))  
  
# Calling this many sites can be very time consuming.  
mam.dl <- get\_download(mam.set)

Sites are assigned to time-period bins as in Graham et al. (1996), Modern (< .5 ka cal BP), Late Holocene (.5 - 4 ka cal BP), Early-Mid Holocene (4 - 10 ka cal BP), Late Glacial (10 - 15 ka cal BP), Full Glacial (15 - 20 ka cal BP) and Late Pleistocene (> 20 ka cal BP). The first step is to build a large table with time and xy coordinates for each site. Time data in "sample.meta" for the mammal data is not the same as for for pollen where age is commonly found in ...$sample.meta$age. Most vertebrate fauna samples, are assigned only younger (sample.meta$age.younger) and older (sample.meta$age.older, Figure 1) bounds with no estimates of mean or median age. Younger and older bounds in this example are averaged, however, this is likely to be methodologically indefensible in the scientific literature. We use it here for illustrative purposes. We use the reshape2 package for R (Wickham 2007) to help manipulate data structures.

compiled.mam <- compile\_downloads(mam.dl)  
  
time.bins <- c(500, 4000, 10000, 15000, 20000)  
  
mean.age <- rowMeans(compiled.mam[, c("age.old", "age.young", "age")], na.rm = TRUE)  
interval <- findInterval(mean.age, time.bins)  
  
periods <- c("Modern", "Late Holocene", "Early-Mid Holocene", "Late Glacial",   
 "Full Glacial", "Late Pleistocene")  
  
compiled.mam$ageInterval <- periods[interval + 1]  
  
mam.melt <- melt(compiled.mam, measure.vars = 10:(ncol(compiled.mam) - 1), na.rm = TRUE,   
 factorsAsStrings = TRUE)  
  
mam.melt$ageInterval <- factor(mam.melt$ageInterval, levels = periods)  
  
mam.lat <- dcast(data = mam.melt, variable ~ ageInterval, value.var = "lat",   
 fun.aggregate = mean, drop = TRUE)[, c(1, 3, 5, 6)]  
  
# We only want taxa that appear at all time periods:  
mam.lat <- mam.lat[rowSums(is.na(mam.lat)) == 0, ]  
  
# Group the samples based on the range & direction (N vs S) of migration. A  
# shift of only 1 degree is considered stationary.  
mam.lat$grouping <- factor(findInterval(mam.lat[, 2] - mam.lat[, 4], c(-11,   
 -1, 1, 20)), labels = c("Southward", "Stationary", "Northward"))  
  
mam.lat.melt <- melt(mam.lat)  
colnames(mam.lat.melt)[2:3] <- c("cluster", "Era")

ggplot(mam.lat.melt, aes(x = Era, y = value)) + geom\_path(aes(group = variable,   
 color = cluster)) + facet\_wrap(~cluster) + scale\_x\_discrete(expand = c(0.1,   
 0)) + ylab("Mean Latitude of Occurance") + theme(axis.text.x = element\_text(angle = 90,   
 hjust = 1))

plot of chunk mammal-example-plot

Even with this fairly simple set of analyses we see that species did not respond uniformly to climatic warming following deglaciation, consistent with the prior work of Graham *et al*. (1996). Although most range shifts were northward, a number of taxa show little change in their ranges and a number show southward range shifts. This example does not examine east-west movement and ignores the issues that may be associated with the complex topography of the mountainous west, or possible confounding effects introduced by temporal variations in the available set of sites. The broader point here is that the use of neotoma can support research that is synchronized with the data holdings of large repositories such as Neotoma and reproducible.

# Conclusion

The whole of the fossil record is much greater than the sum of its parts. Many of our discipline's most important advances were made possible only by the synthesis of many individual fossil occurrences into regional- to global-scale databases of species occurrences, *e.g.*, the Neotoma Paleoecology Database and the Paleobiological Database. Current frontiers in paleoecological informatics include 1) facilitating the input of data into these databases, 2) improved sophistication of the data models employed by these databases, enabling them to handle increasingly complex arrays of paleobiological and associated geochronological data, and 3) enabling the frictionless integration of these resources with other cyberinfrastructure (Uhen et al. 2013; Brewer et al. 2012; Committee 2014).

Here we present the neotoma package for R and show how it can be used to directly transfer data from the Neotoma Paleoecology Database into the R statistical computing environment. The broader goals of this effort are 1) to ease the transfer of data from Neotoma into an environment widely used for paleoecological analyses (Simpson & Oksanen 2014; Simpson 2007; Juggins 2013) and 2) to enable transparent and reproducible scientific workflows. The neotoma package itself is available either from the CRAN repository, or from GitHub (<http://github.com/ropensci/neotoma>) where ongoing open-source development continues. Suggestions for improvement and new code contributions by readers and users are welcome.

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