

# Biomass Prediction All Data (Moving Window) three vars

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This script tiles the study area into  $400m^2$  cells and computes a biomass prediction for each cell using the regression parameters developed in the script All Data Biomass Regression.

```
In [36]: %matplotlib inline
import pandas, numpy, math
from os import system
import time
from matplotlib import pyplot as plt
from IPython.display import display
import matplotlib
import warnings
warnings.filterwarnings('ignore')
from numpy import log as ln
```

Import first and last return textfiles.

```
In [37]: firstfile = r"D:\Users\jkelly\first_all\first_all_combined.txt"
first = pandas.read_table(firstfile, header=0, delimiter=" ")
lastfile = r"D:\Users\jkelly\last_all\last_all_combined.txt"
last = pandas.read_table(lastfile, header=0, delimiter=" ")
system('say table finished')
```

Out[37]: 1

This function computes only last-return mean and first-return 50th percentile cutoff for each grid cell, the explanatory variables identified in the prior script. These are then combined with the regression parameters from the previous script to predict a biomass estimate on the 0.04 ha grid cells.

```
In [46]: def heightstats(L_SERIES, F_SERIES, INT, B1, B2, B3):

    L_MEAN = numpy.mean(L_SERIES)
    H50 = numpy.percentile(F_SERIES, 50)

    BIO = numpy.exp(INT + B1*ln(L_MEAN) + B2*ln(H50) + B3*H50)
    outlist = [L_MEAN, H50, round(BIO,2)]

    return(outlist)
```

The next section sets up the moving window loop. It first builds a pandas dataframe with column names and then sorts on x values in order to work efficiently in left-to-right columns across the study area. It also identifies the extent coordinates and  $\Delta$ , the increment for cell size (here 20x20).

```
In [39]: df = pandas.DataFrame(first)
df.columns = ['Xcoord', 'Ycoord', 'Zcoord', 'Angle']
dfF = df.sort_values(by=['Xcoord'], ascending=True)
dfF.index = range(1, len(dfF) + 1)
df = pandas.DataFrame(last)
df.columns = ['Xcoord', 'Ycoord', 'Zcoord', 'Angle']
dfL = df.sort_values(by=['Xcoord'], ascending=True)
dfL.index = range(1, len(dfL) + 1)
print("dfF = " + str(len(dfF)))
print("dfL = " + str(len(dfL)))
xmin0 = 604790.00
xmax0 = 610380.37
delta = 20
ymin0 = 6596136.30
ymax0 = 6600916.67
```

```
dfF = 188881266
```

```
dfL = 41918199
```

```
In [40]: dfL.head()
```

```
Out[40]:
```

	Xcoord	Ycoord	Zcoord	Angle
1	604924.18	6598334.93	0.01	3
2	604924.20	6598335.88	0.01	3
3	604924.24	6598332.00	0.01	3
4	604924.27	6598329.08	0.01	3
5	604924.29	6598328.11	0.01	3

The next section uses  $\Delta$  to create lists of x and y block endpoint coordinates.

```
In [41]: ylist = []
# from max to min by 20s
for i in range(int(ymin0), int(math.ceil(ymax0)), int(delta)):
    ylist.append(i)
xlist = []
for i in range(int(xmin0), int(math.ceil(xmax0)), int(delta)):
    xlist.append(i)

len(xlist), len(ylist)
```

```
Out[41]: (280, 240)
```

Dimensions arrays to hold statistics.

```
In [65]: f = open('parameters_all.csv', 'r')
        with f:
            intercept = float(f.readline())
            b1 = float(f.readline())
            b2 = float(f.readline())
            b3 = float(f.readline())

        f.close()
        print (intercept, b1, b2, b3)

0.691575840756 0.134929275669 -0.318404481456 0.128437207872
```

The next section loops over x values. It isolates a column by slicing off a  $\Delta$ -unit-wide xstrip, which it then sorts by y values in order to work in blocks from bottom to top, chopping off each  $\Delta$ -unit-wide yblock and calculating that block's lidar metrics and predicted biomass.

```
In [47]: statslist=[] ## Good to here

        for xmin in xlist:
            xmax = xmin + delta

            # cut x-strip between current xmin and xmax
            xstripF = dfF.loc[lambd df2:(df2.Xcoord >= xmin)
                             & (df2.Xcoord < xmax), :]
            xstripL = dfL.loc[lambd df2:(df2.Xcoord >= xmin)
                             & (df2.Xcoord < xmax), :]

            if len(xstripF) == 0: continue
            #print("*** xstrip centered at " + str(xmin + (delta/2)) +
            #      " has " + str(len(xstrip)) + " elements")

            # sort x-strip by y
            xstripF = xstripF.sort_values(by=['Ycoord'], ascending=True)
            xstripF.index = range(1, len(xstripF) + 1)
            xstripL = xstripL.sort_values(by=['Ycoord'], ascending=True)
            xstripL.index = range(1, len(xstripL) + 1)

            for ymin in ylist:
                ymax = ymin + delta

                # cut y-block between current ymin and ymax
                yblockF = xstripF.loc[lambd df3:(df3.Ycoord >= ymin)
                                       & (df3.Ycoord < ymax), :]
                yblockL = xstripL.loc[lambd df3:(df3.Ycoord >= ymin)
                                       & (df3.Ycoord < ymax), :]

                if len(yblockF) < 10: continue
                elif len(yblockF) >= 10:
```

```

blocklist = heightstats(yblockL.Zcoord, yblockF.Zcoord, interco
anglemean = numpy.mean(yblockF.Angle)
blocklist.append(anglemean)
blocklist.append(xmin + delta/2)  # append block center coords
blocklist.append(ymin + delta/2)
statslist.append(blocklist)

```

```

system('say process finished') #only works on mac

```

Out[47]: 1

Converts statistics list to pandas dataframe and writes to csv.

```

In [61]: #statsarray = numpy.array(statslist)
outdf = pandas.DataFrame(statslist)
columnlist = ['L_mean', 'F_h50', 'Biomass_est', 'Angle',
              "Xcenter", "Ycenter"]
outdf.columns = columnlist
outdf.Biomass_est = outdf.Biomass_est.apply(pandas.to_numeric)
outdf.to_csv('trial_run_1.csv')

```

In [67]: outdf[0:100]

```

Out[67]:
   L_mean  F_h50  Biomass_est  Angle  Xcenter  Ycenter
0  1.708929  12.940      5.01  20.000000  604920.0  6597886.0
1  1.649508   7.530      2.95  19.074380  604920.0  6597906.0
2  2.048000   9.000      3.47  18.620690  604920.0  6597926.0
3  0.863462   8.250      2.88  17.989247  604920.0  6597946.0
4  0.215000   8.310      2.40  17.020833  604920.0  6597966.0
5  0.625429   7.545      2.60  16.352273  604920.0  6597986.0
6  0.493529   2.960      1.88  15.720430  604920.0  6598006.0
7  0.572083   5.050      2.12  14.961905  604920.0  6598026.0
8  0.240000   5.430      1.93  14.000000  604920.0  6598046.0
9  0.890000   6.260      2.45  13.341772  604920.0  6598066.0
10 1.175769  11.845      4.25  12.702128  604920.0  6598086.0
11 3.719259  17.610      9.18  12.000000  604920.0  6598106.0
12 7.320303  17.050      9.46  11.053097  604920.0  6598126.0
13 4.144545  16.060      7.86  10.269231  604920.0  6598146.0
14 7.368462  18.130     10.67   9.600000  604920.0  6598166.0
15 4.234231  20.090     12.32   8.895652  604920.0  6598186.0
16 5.710909  19.025     11.38   8.000000  604920.0  6598206.0
17 7.990588  18.020     10.65   7.161017  604920.0  6598226.0
18 5.662917  15.765      7.94   6.434426  604920.0  6598246.0
19 4.478718  17.300      9.10   5.702479  604920.0  6598266.0
20 1.459487  10.630      3.88   4.957265  604920.0  6598286.0
21 4.641591  15.190      7.27   4.000000  604920.0  6598306.0
22 6.033276  16.110      8.32   3.211679  604920.0  6598326.0

```

23	4.535833	0.010	10.62	2.510204	604920.0	6598346.0
24	NaN	0.010	NaN	1.783784	604920.0	6598366.0
25	NaN	0.010	NaN	0.987013	604920.0	6598386.0
26	NaN	0.010	NaN	0.000000	604920.0	6598406.0
27	NaN	0.010	NaN	0.000000	604920.0	6598426.0
28	NaN	0.010	NaN	-0.423077	604920.0	6598446.0
29	NaN	0.010	NaN	-1.170455	604920.0	6598466.0
..	...	...	...	...	...	...
70	6.472672	19.605	12.36	8.000000	604940.0	6598206.0
71	7.279588	20.190	13.41	7.192067	604940.0	6598226.0
72	5.837090	18.950	11.32	6.443320	604940.0	6598246.0
73	5.920509	17.180	9.32	5.714579	604940.0	6598266.0
74	4.432748	16.090	7.96	4.974257	604940.0	6598286.0
75	5.549938	15.900	8.04	4.002165	604940.0	6598306.0
76	4.851095	15.500	7.56	3.263830	604940.0	6598326.0
77	4.390476	1.880	2.54	2.536082	604940.0	6598346.0
78	0.010000	0.010	4.65	1.762082	604940.0	6598366.0
79	NaN	0.010	NaN	1.000000	604940.0	6598386.0
80	NaN	0.010	NaN	0.012658	604940.0	6598406.0
81	NaN	0.010	NaN	0.000000	604940.0	6598426.0
82	NaN	0.010	NaN	-0.405063	604940.0	6598446.0
83	NaN	0.010	NaN	-1.160643	604940.0	6598466.0
84	NaN	0.010	NaN	-2.000000	604940.0	6598486.0
85	NaN	0.010	NaN	-2.831933	604940.0	6598506.0
86	4.967000	0.960	2.84	-3.694190	604940.0	6598526.0
87	3.478871	7.840	3.36	-4.286501	604940.0	6598546.0
88	1.261613	0.190	3.58	-5.000000	604940.0	6598566.0
89	0.230000	0.080	3.70	-5.981550	604940.0	6598586.0
90	0.110000	0.010	6.43	-6.668142	604940.0	6598606.0
91	NaN	0.010	NaN	-7.365297	604940.0	6598626.0
92	NaN	0.010	NaN	-8.020942	604940.0	6598646.0
93	NaN	0.010	NaN	-9.000000	604940.0	6598666.0
94	NaN	0.010	NaN	-9.729084	604940.0	6598686.0
95	NaN	0.010	NaN	-10.376471	604940.0	6598706.0
96	NaN	0.010	NaN	-11.058559	604940.0	6598726.0
97	NaN	0.080	NaN	-12.000000	604940.0	6598746.0
98	NaN	0.010	NaN	-12.661905	604940.0	6598766.0
99	NaN	0.010	NaN	-13.364583	604940.0	6598786.0

[100 rows x 6 columns]

```
In [68]: outdf1 = outdf.replace([numpy.inf, -numpy.inf],
                                numpy.nan).dropna(subset=['Biomass_est'], how="all")
```

These are the total biomass estimate and the overall biomass-per-hectare for the 2170.5 hectare study area:

```
In [69]: total = round(outdf1['Biomass_est'].sum(), 3)
```

```
In [70]: print ("total biomass in Mg, average biomass per hectare")
         "{:,}".format(total), "{:,}".format(round(total/2170.5, 3))
```

total biomass in Mg, average biomass per hectare

```
Out[70]: ('212,177.71', '97.755')
```

```
In [71]: dfL.mean()
```

```
Out[71]: Xcoord      6.075622e+05
         Ycoord      6.598420e+06
         Zcoord      2.485655e+00
         Angle       1.475276e+00
         dtype: float64
```

```
In [72]: dfF.mean()
```

```
Out[72]: Xcoord      6.076169e+05
         Ycoord      6.598451e+06
         Zcoord      8.214472e+00
         Angle       1.024749e+00
         dtype: float64
```

```
In [73]: outdf1.mean()
```

```
Out[73]: L_mean      2.162664e+00
         F_h50      8.501369e+00
         Biomass_est  4.261966e+00
         Angle       1.756524e+00
         Xcenter     6.075992e+05
         Ycenter     6.598487e+06
         dtype: float64
```

```
In [ ]:
```

```
In [ ]:
```