Impacts of Toxic Pollution on Life Expectancy and Cancer Rates for Local Communities

Introduction to Machine Learning, Final Project Report

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Where applicable, I've labelled the sections with indices from the Project requirements as found in doc/Project Requirements.pdf for reference. Please do read the full document however as all sections are still relevant.

Problem Overview (3.a)

Using publicly available data from the EPA and CDC, I would like to explore the potential links of toxic dumping to general life expectancy and cancer rates at the local level, throughout the United States. Specifically, I have the data to do so from 2000 to 2016. Initially, I will constrain this range to 2010 to 2016 to align with the most recent Census (needed for life expectancy).

It is well known that certain chemicals are a contributing cause to a decline in public health. However, the process of disposing or sequestering these toxics are not equal, and local ecological factors may mean one method is significantly more dangerous for a given region than other methods.

Motivation (3.b)

Environmental Accountability has historically been lower than required for long term ecological stability. Over time, the lack of accountability compounds to a large implicit cost for communities (and countries as a whole) as soil, air, and water supplies become more saturated with dangerous materials. The ability to identify a correlation between toxic dumping and public health is an important first step in fixing affected communities.

Therefore, my goal is to identify counties within the United States which have an above-average amount of industrial waste processing, and assess that area's life expectancy or cancer rates against national averages.

Methodology (5.a)

I have separated my data into three categories: Health (cancer and life expectancy), TRI (the Toxic Release Inventory), and the AQI (Air Quality Index). Given these three categories, I will determine some geo-spatial correlation between their values. Data is grouped by various regions including state, county, and custom defined regions. I will use OpenStreetMap's Nominatim API to convert these regions to their Latitude and Longitude. State level information will be folded into county-level data points, and custom regions will be handled by changing the precision of the Lat/Lon measurement to integer accuracy, grouping neighboring regions into one. This represents areas roughly 100km wide.

References: Data Overview (4.)

Four datasets are considered for this project from 2010 to 2016. For measures of Air Quality, I consider the EPA's AirData AOI (Air quality index) data set. For measures of toxicity. I consider the EPA's Toxics

```
In [1]: ## Work in the root directory
        %cd
        /home/cole/Desktop/intro-to-ml-final-project
In [2]: import json
        import geocoder
        from sklearn.model_selection import train_test_split
        from sklearn preprocessing import MaxAbsScaler
In [3]: def make_lat_lon_map(inputs, output_json_file='data/latlon.json', load_from='data/
            Queries OpenStreetMap for Lat/Lon pairs given a list
            of input queries. The results are incrementally written
            to an output file in `data` by default. Errors are logged
            to a separate file for manual lookup.
                                        A list of UNIQUE values to query
            :param inputs:
            :param output_json_file:
                                        Optional separate JSON output file
                                        Resumes from another JSON file
            :param load from:
                                        entries in the file will not be
                                        reprocessed
            :return: `None`
            # {"query": [lat, lon]}
            json_acc = {}
            # Remove duplicates
            inputs = list(set(inputs))
            if load from:
                with open(load_from, 'r') as json_file:
                    json_acc = json.load(json_file)
                inputs = [i for i in inputs if i not in json acc.keys()]
            def dump(j):
                with open(output json file, 'w') as fp:
                    json.dump(j, fp, indent=4)
            def dump_errors(e):
                with open("data/nominatim_errors.txt", 'w') as fp:
                    for err in e:
                        fp.write(err + '\n')
            errors = []
            for query in inputs:
                query = str(query)
                if query not in json acc:
                    g = geocoder.osm(query)
                    if not q.ok:
                        print("[ WRN ]", query, "--", str(g))
                        errors.append(query)
                        dump errors(errors)
                    else:
                        print("[ MAP ] Found", g.latlng, "for", query)
                        json acc[query] = g.latlng
                        dumn(ison acc)
```

Common Loading (Originally in src/etl/load.py)

We will first define a common loading procedure that works for most of the datasets considered here.

```
In [4]: import glob
import json
import os
import re
import nandas as nd
```

```
In [5]: # Not required for Notebook, function copied in
        # from src.etl.preprocess import make_lat_lon_map
        RAW DATA = 'data/raw/'
        PREPARED_DATA = 'data/prepared'
        # Load the Lat/Lon Informtion
        with open('data/latlon.json', 'r') as fp:
            lat_lon_json = json.load(fp)
        def _load(subpath):
            Common loadining method for CSVs in data directory.
            :param subpath:
                                Subpath in `data`
            :return:
                                DataFrame
            path = os.path.join(RAW_DATA, subpath, '*.csv')
            all_files = glob.glob(path)
            li = []
            for filename in all_files:
                df = pd.read_csv(filename, index_col=None)
                li.append(df)
            return pd.concat(li, axis=0, ignore_index=True)
        def _get_lat_lon(q):
            Returns latitude and longitude pair from generated
            JSON file.
            :param q:
                        String query
            :return:
                        Latitude Longitude Pair
            if q in lat lon json:
                return lat lon json[q]
            else:
                return [None, None]
        def _change_precision(a, prec=0):
            Changes the precision of the Latitude
            or Longitude arrays while preserving
            sparsity.
                            Input array
            :param a:
            :param prec: New Decimal Precision
            :return:
                            Rounded array
            acc = []
            for e in a:
                if e is None:
                    acc.append(None)
                    acc.append(round(float(e), prec))
            return acc
```

TRI

We will begin by preparing the TRI dataset. In order to get this into a common format, we will make the following transformations:

- 1. Clean the column names to remove spaces and capitals
- 2. Select relevant columns (as defined in function)
- 3. Pivot on industry since it is categorical
- 4. Fix the schema for the columns which will become the index (lat, lon, year)
- 5. Change precision of lat and lon
- 6. Regroup by year, lat, lon and aggregate by sum

The TRI dataset is fortunate enough to not require lat, lon conversion. It comes with lat lon already.

```
In [6]: | def load_tri():
                 Loads the TRI CSV files into one DataFame
                  :return: DataFrame
                 # Remove NA columns
                 combined = _load('tri')
                 # Clean column names
                 cols = combined.columns
                 cols = [c.split(' - ')[-1] for c in cols]
cols = [c.split('. ')[-1].lower() for c in cols]
                 combined.columns = cols
                 combined = combined.rename(columns={'production wste (8.1-8.7)': 'production v
                                                                       'latitude': 'lat', 'longitude': 'lon'})
                 # Select Relevant
                 'underground cl i', 'underground c ii-v', 'landfills',
'rcra c landfill', 'other landfills', 'land treatment',
'surface impndmnt', 'rcra surface im', 'other surface i',
'on-site release total', 'trns rlse', 'trns trt',
'total transfers', 'm10', 'm41', 'm62', 'm40 metal', 'm61 metal', 'm71'
'm81', 'm82', 'm72', 'm63', 'm66', 'm67', 'm64', 'm65', 'm73', 'm79',
'm90', 'm94', 'm99', 'off-site release total', 'm20', 'm24', 'm26',
'm28', 'm93', 'off-site recycled total', 'm56', 'm92',
'm40 non-metal', 'm50', 'm54'
                            'm40 non-metal', 'm50', 'm54',
'm61 non-metal', 'm69', 'm95', 'off-site treated total',
'total transfer', 'total releases', 'releases', 'on-site contained',
                            'off-site contain', 'production waste']
                 combined = combined[sel].dropna(axis=1)
                 combined.columns = [c.replace(' ', '_') for c in combined.columns]
                 # Pivot industry
                 d = pd.get dummies(combined['industry sector'], prefix='sector')
                 df = pd.concat([combined, d], axis=1).drop(['industry sector'], axis=1)
                 # Fix schema
                 df = df.dropna()
                 df.year = df.year.astype(int)
                 df.lat = df.lat.astype(int)
                 df.lon = df.lon.astype(int)
                 # Shift Precision
                 df.lat = change precision(df.lat)
                 df.lon = _change_precision(df.lon)
                 df = df.dropna().groupby(by=['year', 'lat', 'lon'], as index=False).sum()
                  return df dronna()
```

AQI

We will begin by preparing the AQI dataset. In order to get this into a common format, we will make the following transformations:

- 1. Clean the column names to remove spaces and capitals
- 2. Select relevant columns (as defined in function)
- 3. Convert date to year by remove the month and day
- 4. Pivot on defining_parameter since it is categorical. This column denotes the worst factor for air quality that day
- 5. Regroup by 'year', 'state_name', 'county_name', aggregating by mean.
- 6. Convert 'state_name', 'county_name' to lat and lon using the JSON file.
- 7. Fix the schema for the columns which will become the index (lat, lon, year)
- 8. Change precision of lat and lon
- 9. Regroup by year, lat, lon and aggregate by ${\bf sum}$

```
In [7]: def load_aqi():
            Loads the AQI CSV files into one DataFame
            :return: DataFrame
            # Remove NA columns
            combined = _load('aqi').dropna(axis=1)
            # Select relevant columns
            sel = ['State Name', 'county Name', 'Date', 'AQI', 'Defining Parameter']
            combined = combined[sel]
            # Clean column names
            combined.columns = [c.replace(" ", "_").lower() for c in combined.columns]
            # Convert date column to year only
            combined.date = combined.date.str.split('-').str[0]
            combined = combined.rename(columns={'date': 'year'})
            combined = combined.astype({'year': 'int64'})
            # Pivot defining_parameter column
            d = pd.get_dummies(combined['defining_parameter'], prefix='defining')
            df = pd.concat([combined, d], axis=1).drop(['defining_parameter'], axis=1)
            df = df.dropna().groupby(by=['year', 'state_name', 'county_name'], as_index=F@index
            # Convert to Lat/Lon
            loc = list(df.state_name + ", " + df.county_name)
            make lat lon map(list(set(loc)))
            lat = [_get_lat_lon(i)[0] for i in loc]
            lon = [_get_lat_lon(i)[1] for i in loc]
            df['lat'] = _change_precision(lat)
            df['lon'] = change precision(lon)
            # Fix Schema
            df = df.dropna()
            df.year = df.year.astype(int)
            df.lat = df.lat.astype(int)
            df.lon = df.lon.astype(int)
            # Group up
            df = df.dropna().groupby(by=['year', 'lat', 'lon'], as_index=False).sum()
            return df dronna()
```

Cancer

We will begin by preparing the Cancer dataset. In order to get this into a common format, we will make the following transformations:

- 1. Clean the column names to remove spaces and capitals
- 2. Select relevant columns (as defined in function)
- 3. Pivot on leading_cancer_sites since it is categorical.
- 4. Fix the MSA column by defining a custom transformation policy for use in querying with **Nominatum**. This is required since the MSA regions sometimes include metropolitan area spanning across state lines for which Nominatim isn't able to handle.
- 5. Convert the cleaned MSA column to lat and lon.
- 6. Fix the schema for the columns which will become the index (lat, lon, year)
- 7. Change precision of lat and lon
- 8. Regroup by year, lat, lon and aggregate by sum

```
In [8]: def load cancer():
            path = os.path.join(RAW_DATA, "health/United States and Puerto Rico Cancer Sta
            cancer = pd.read_csv(path, delimiter='\t')
            # Select relevant columns
            sel = ['Year', 'Leading Cancer Sites', 'MSA', 'Count', 'Population', 'Age-Adj
            cancer = cancer[sel]
            cancer.columns = [c.replace(' ', '_').lower() for c in cancer.columns]
            # Pivot criteria
            d = pd.get dummies(cancer['leading cancer sites'], prefix='cancer')
            d = d.multiply(cancer['count'], axis="index")
            cancer = pd.concat([cancer, d], axis=1).drop(['leading cancer sites', 'count']
            # Fix MSA
            tmp = cancer.msa
            ntmp = []
            # Also load manual map
            manual = \{\}
            with open('data/cancer_manual_map.txt', 'r') as fp:
                 for l in fp:
                     items = l.split('|')
                     manual[items[0].strip()] = items[1].strip()
            # MSA Conversion Policy
            for t in tmp:
                t = str(t)
                t = t.replace('-', '')
if len(re.split(',|,,', t)) == 1:
                     ntmp.append(t)
                     continue
                 states = re.split(',|,,', t)[1]
                 state = states.split()[0]
                 areas = t.split()
                 if len(areas) > 1:
                     area = t.split()[0] + ' ' + t.split()[1]
                 else:
                    area = areas[0]
                tt = area + ", " + state
                 if tt in manual:
                    ntmp.append(manual[tt])
                 else:
                     ntmp.append(tt)
            cancer.msa = ntmp
            cancer = cancer.dropna()
            # Convert to Lat/Lon
            loc = list(cancer.msa)
            make lat lon map(list(set(loc)))
            lat = [qet lat lon(i)[0] for i in loc]
            lon = [\_get\_lat\_lon(i)[1]  for i  in loc]
            cancer['lat'] = change precision(lat, prec=0)
            cancer['lon'] = _change_precision(lon, prec=0)
            # Fix Schema
```

Life Expectancy

We will begin by preparing the Life Expectancy dataset. In order to get this into a common format, we will make the following transformations:

- 1. Clean the column names to remove spaces and capitals
- 2. Select relevant columns (as defined in function)
- 3. Fix the Life Exp range to Min and Max values, which could potentially be good features.
- 4. Group by state and county, aggregate by mean to get the average life exp for that state and county
- 5. Convert state and county to lat and lon using the JSON file
- 6. Fix the schema for the columns which will become the index (lat, lon, year)
- 7. Change precision of lat and lon
- 8. Regroup by year, lat, lon and aggregate by sum

```
In [9]: def load life exp():
            path = os.path.join(RAW_DATA, 'health', 'U.S._Life_Expectancy_at_Birth_by_Stat
            life = pd.read_csv(path)
            # Select Relevant
            sel = ['State', 'County', 'Life Expectancy',
                    'Life Expectancy Range']
            life = life[sel].dropna()
            # Fix column names
            life.columns = [c.replace(' ', ' ').lower() for c in life.columns]
            # Fix Life Exp Range to Min and Max
            rng = list(life.life expectancy range)
            rng max = [float(str(r).split('-')[1].strip()) for r in rng]
            rng min = [float(str(r).split('-')[0].strip()) for r in rng]
            life['life_expectancy_max'] = rng_max
            life['life_expectancy_min'] = rng_min
            life = life.rename(columns={'life_expectancy': 'life_expectancy_avg'})
            life = life.drop('life_expectancy_range', axis=1).dropna()
            life = life.groupby(['state', 'county'], as_index=False).mean()
            # Convert to Lat/Lon
            print("Getting lat lon list")
            loc = list(life.state + ", " + life.county)
            make_lat_lon_map(list(set(loc)))
            lat = [_get_lat_lon(i)[0] for i in loc]
            lon = [_get_lat_lon(i)[1] for i in loc]
            life['lat'] = _change_precision(lat, prec=0)
            life['lon'] = _change_precision(lon, prec=0)
            # Fix Schema
            life = life.dropna()
            life.lat = life.lat.astype(int)
            life.lon = life.lon.astype(int)
            life = life.dropna().groupby(by=['lat', 'lon'], as index=False).sum()
            return life.dropna()
```

Load All

Combine the datasets by year, lat, and lon. This is then written to a file that will constitute our main dataset.

```
In [10]: def load_all(from_file=True):
              if from_file:
                   return pd.read csv('data/merged.csv')
              print("[ LOAD ] Loading AQI...")
              aqi = load aqi()
              print(aqi)
              print("[ LOAD ] Loading TRI...")
              tri = load_tri()
              print(tri)
              print("[ LOAD ] Loading Life...")
              life = load_life_exp()
              print(life)
              print("[ LOAD ] Loading Cancer...")
              cancer = load_cancer()
              print(cancer)
              # Join
              j = aqi.merge(tri, on=['year', 'lat', 'lon'], how='inner')
j = j.merge(cancer, on=['year', 'lat', 'lon'], how='inner')
              j = j.merge(life, on=['lat', 'lon'], how='left')
              j = j.dropna()
              # TODO: Fix weird behavior with duplicated rows in merge that make this line i
              j = j.groupby(by=['year', 'lat', 'lon'], as index=False).mean()
              j.to_csv('data/merged.csv', index=False)
              return i
```

```
In [11]: df = load_all(from_file=True)
Out[11]:
```

	year	lat	lon	aqi	defining_CO	defining_NO2	defining_Ozone	defining_PM10	defining_PM2.5 (
0	2010	21	-158	39.715068	0.000000	0.00274	0.230137	0.024658	0.742466
1	2010	26	-98	44.077348	0.000000	0.00000	0.406077	0.002762	0.591160
2	2010	27	-82	42.367123	0.000000	0.00000	0.695890	0.010959	0.293151
3	2010	28	-82	183.767123	0.000000	0.00000	3.038356	0.002740	0.821918
4	2010	28	-81	208.062127	0.000000	0.00000	3.714128	0.000000	1.285872
					•••	•••			
581	2016	43	-74	104.521436	0.000000	0.00000	1.748634	0.000000	1.251366
582	2016	44	-116	34.267218	0.000000	0.00000	0.000000	0.000000	1.000000
583	2016	45	-123	30.764543	0.000000	0.00000	0.407202	0.000000	0.592798
584	2016	47	-122	103.114754	0.000000	0.31694	0.989071	0.000000	1.693989
585	2016	48	-117	85.160502	0.002732	0.00000	0.352459	0.169399	2.475410

586 rows × 125 columns

Data Evaluation: High Level Statistics

Let us view some high level statistics about our dataset. Namely, we should look at the:

- 1. Number of Rows
- 2. Number of Columns
- 3. Columns with Highest Variance
- 4. Columns with Lowest Variance
- 5. Some basic correlations

```
In [12]: |print("Number of Rows:", len(df))
         print("Number of Cols:", len(df.columns), '\n')
         print("Highest Variance Cols:")
         print(df.var().sort_values(ascending=False)[:10], '\n')
         print("Lowest Variance Cols:")
         print(df.var().sort values()[:10], '\n')
         print("Top 20 Correlations to Age-Adjusted Rate:")
         print(df[[c for c in df.columns if 'cancer' not in c]]
               _corr()['age-adjusted_rate'] sort_values(ascending=False)[:201)
         Number of Rows: 586
         Number of Cols: 125
         Highest Variance Cols:
         population
                                    3.779196e+15
         production waste
                                    2.542578e+11
         total releases
                                    1.909209e+10
         on-site release total
                                   1.894071e+10
         other surface i
                                    5.614605e+09
         total transfer
                                    1.802190e+09
         off-site_recycled_total
                                   1.550502e+09
         m24
                                    1.442020e+09
                                    2.049389e+08
         water
         off-site_release_total
                                    1.271601e+08
         dtype: float64
         Lowest Variance Cols:
         releases
                               0.000000e+00
         landfills
                               0.000000e+00
         underground
                               0.000000e+00
         m63
                               0.000000e+00
         m71
                               0.000000e+00
                               0.000000e+00
         surface_impndmnt
         sector_Publishing
                               3.180443e-07
         sector_Apparel
                               6.450643e-07
         sector_Coal Mining
                               2.369974e-05
         sector_Tobacco
                               5.922906e-05
         dtype: float64
         Top 20 Correlations to Age-Adjusted Rate:
         age-adjusted rate
                                               1.000000
         population
                                               0.495891
                                               0.422032
         defining_PM2.5
                                               0.419695
         aqi
         defining_Ozone
                                               0.390418
         defining NO2
                                               0.371733
         sector_Petroleum Bulk Terminals
                                               0.357192
                                               0.314356
         life expectancy min
                                               0.314105
                                               0.305783
         life expectancy max
         life_expectancy_avg
                                               0.291387
         sector_Tobacco
                                               0.286238
         sector Textiles
                                               0.215277
         sector_Miscellaneous Manufacturing
                                               0.192700
         m93
                                               0.160954
         m95
                                               0.139101
         defining CO
                                               0.137515
         underground cl i
                                               0.136676
                                               0.116186
         sector Electric Utilities
                                               0.104435
         Name: ane-adjusted rate dtyne: float61
```

Interpretation

From above, we can see that there are no columns in our desgin matrix that are immediately strongly correlated with the Age-Adjusted Rate of cancer levels for a given year and area.

This was expected however. Efforts to relate toxic dumping with cancer rates has been historicall inconclusive as evidenced by the following paper survey.

Hazardous waste and health impact: a systematic review of the scientific literature

https://www.ncbi.nlm.nih.gov/pmc/articles/PMC5637250/ (https://www.ncbi.nlm.nih.gov/pmc/articles/PMC5637250/ (https://www /PMC5637250/)

In this survey study, the authors assessed 57 papers on the health impacts of toxic waste. The overall results from many studies were inconclusive - There are many confounding factors which contribute to cancer rates among individuals. However, they found that oil industry waste to be the most hazardous of the types considered in the survey. This study is fairly recent, published in 2017, but again considers a traditional sample of selected communities, instead of a national survey.

Experiment: Visualization (Originally src/model /cluster.py)

We can now view the prepared datasets individally and together to assess potential clusters, and to help select our model later on.

We will use a 2-component PCA Decomposition on the fully prepared dataset. Let us define a generalized visualization function.

In [13]: from matplotlib import pyplot as plt from sklearn decomposition import PCA

```
In [14]: def view pca(df, emphasize column=None, title=None, xlabel=None, ylabel=None):
             Charts the 2-component PCA decomposition of
             the given DataFrame, with optional
             coloring upon a separate column. This
             column will NOT be part of the PCA decomposition.
             :param df:
                                         DataFrame to decompose
             :param emphasize_column:
                                         Column for coloring PCA
             :param title:
                                         Title of Plot
             :param xlabel:
                                         X axis label
             :param ylabel:
                                         Y axis label
                                         `None`
             :return:
             c = None
             if emphasize column:
                 c = df[emphasize column]
                 df = df.drop(emphasize column, 1)
             pca = PCA().fit_transform(df)
             plt.scatter(pca[:, 0], pca[:, 1], c=c)
             plt.title(title)
             plt.xlabel(xlabel)
             plt.ylabel(ylabel)
             nlt show()
```

Defining our Y columns

In order to have a meaningful clustering example, we need to define our Y columns and remove them from the DataFrame before the PCA decomposition. The following function will **officially define the columns of our Y set**.

```
In [15]: def _in_terms(t):
    y_terms = ['age-adjusted_rate', 'count', 'cancer', 'population', 'life_expectation'
    for ty in y_terms:
        if ty in t:
            return True
    return False

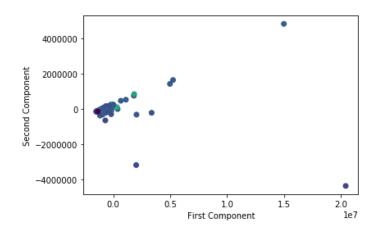
y_cols = [c for c in df.columns if _in_terms(c)]
x_cols = [c for c in df.columns if c not in y cols]
```

All Merged Data w. EMA, Colored by Age-Adjusted Cancer Rates

Visualize the initial clusters of age-adjusted cancer rates for each area. We will group each area by adding up all the years observed.

```
In [16]: # Sum over all years
grouped = df.groupby(by=['lat', 'lon']).sum()
drop cols = [v for v in v cols if v != 'age-adjusted rate']
```

All Merged Data, Colored by Age-Adjusted Cancer Rates

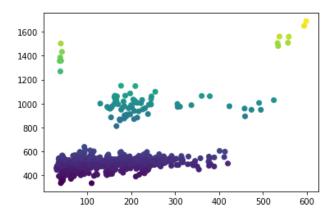


We can immediately see some outliers, but let's take a look at some specific featuers as they relate to the age adjusted rate

Age Adjusted Rate vs AQI (Colored: Age Adjusted Rate)

In [18]: nlt scatter(df['ani'] df['ane-adjusted rate'] c=df['ane-adjusted rate'])

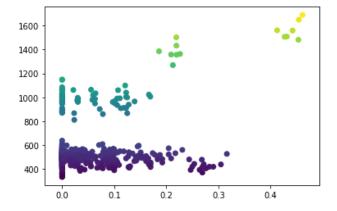
Out[18]: <matplotlib.collections.PathCollection at 0x7fa3cc06c510>



This looks really good. We can tell the realtionship isn't linear, but there does exist defined clusters which pcoket nicely with the age adjusted rate!

Age Adjusted Rate vs Number of Dumpings by Petroleum Companies (Colored: Age Adjusted Rate)

In [19]: | nlt_scatter(df['sector Petroleum Bulk Terminals'] | df['age_adjusted_rate'] | c=df[]
Out[19]: | <matplotlib.collections.PathCollection at 0x7fa3c7b32690>



This one is also pretty good. We notice again that there are clusters of high-count of dumping from Petroleum Companies, with high cancer rates as well.

Algorithm: Feature Engineering (Originally src/model/features.py)

After our data is loaded into a single Pandas DataFrame, we now engineer some features that might be helpful. One such feature is the **exponential moving average** for each feature column. I do this to model the time-weighted importance of each factor which may contribute to life expectancy and cancer rates, with the assumption that **areas will consistently high levels of air pollution and toxic dumping exhibit higher rates and lower life expectancy than areas which have a higher recent spike in air pollution and toxic dumping**. The EMA will capture this sentiment using one, two, and three year lags.

Each feature will become a separate EMA column, denoted by _emaX where X is either 1, 2, or 3.

```
In [20]: def sma_featurizer(df, by=None, lags=None, ema=True):
              Computes the SMA of each column, grouped by
             the `by` set. This follows for each `i` in `lags`. The default `by` is `[lat, lon]`. The default `lags` is `[1,2,3]`, which represent years
              in the merged dataset in this project. SMA
              columns will be appended with prefix smaN .
                              DataFrame to featurize
              :param df:
              :param by: Grouping columns (default: [lat, lon]
              :param lags: SMA i periods (default: [1,2,3])
              :param ema: Use exponential moving averages
              :return:
                            Featurized DataFrame
              # Resolve defaults
              if by is None:
                  by = ['lat', 'lon']
              if lags is None:
                  lags = [1, 2, 3]
              # Re-index DataFrame to by columns plus year
             df = df.set_index(by + ['year'])
              # Group by without year to compute SMA/EMA
              g = df.groupby(by=by)
              # Define an inner convenience function
              def roll(i):
                  # Apply the SMA/EMA to the inner groupings, which should
                  # contain "year"
                  if ema:
                      r = g.apply(lambda x: x.sort_values(by='year').ewm(i).mean())
                      r.columns = ["ema" + str(i) + "_" + c for c in r.columns]
                      r = g.apply(lambda x: x.sort values(by='year').expanding(i).mean())
                      r.columns = ["sma" + str(i) + "_" + c for c in r.columns]
                  return r
              # Compute for all lags
              features = roll(lags[0])
              for lag in lags[1:]:
                  tmp = roll(lag)
                  # Join on by plus year, since year is also still in index
                  features = features.join(tmp, how='inner')
              # Join on the original DataFrame, sort, return
              return of inin(features how='inner') sort index()
```

Featurize our DataFrame

```
In [21]: df_ema = sma_featurizer(df, ema=True).dropna()
df_ema
```

Out[21]:

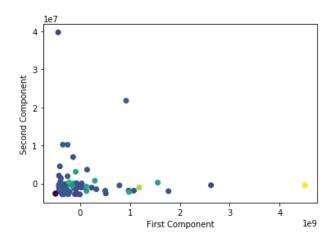
aqi defining_CO defining_NO2 defining_Ozone defining_PM10 defining_PM2.5 defini

lat	lon	year							
21	-158	2010	39.715068	0.000000	0.002740	0.230137	0.024658	0.742466	0
		2011	39.260274	0.000000	0.000000	0.356164	0.016438	0.624658	0
		2012	34.628415	0.000000	0.000000	0.352459	0.013661	0.628415	0
		2013	32.632877	0.000000	0.000000	0.386301	0.008219	0.605479	0
		2014	29.901961	0.000000	0.002801	0.641457	0.044818	0.299720	0
			•••		•••				
48	-117	2012	51.567907	0.005464	0.000000	0.341530	1.010929	0.642077	0
		2013	56.944754	0.008219	0.000000	0.372603	1.024658	0.594521	0
		2014	49.402620	0.016438	0.000000	0.326027	1.147945	0.509589	0
		2015	109.913899	0.000000	0.000000	0.349862	0.414061	2.236077	0
		2016	85.160502	0.002732	0.000000	0.352459	0.169399	2.475410	0

586 rows × 488 columns

```
In [22]: print("Number of Rows:", len(df_ema))
         print("Number of Cols:", len(df_ema.columns), '\n')
         print("Highest Variance Cols:")
         print(df_ema.var().sort_values(ascending=False)[:10], '\n')
         print("Lowest Variance Cols:")
         print(df ema.var().sort values()[:10], '\n')
         print("Top 20 Correlations to Age-Adjusted Rate:")
         print(df ema[[c for c in df ema.columns if 'cancer' not in c]]
               .corr()['age-adjusted_rate'].sort_values(ascending=False)[:201)
         Number of Rows: 586
         Number of Cols: 488
         Highest Variance Cols:
         population
                                 3.779196e+15
         emal population
                                 3.742045e+15
         ema2 population
                                 3.724200e+15
         ema3 population
                                 3.714826e+15
         production waste
                                 2.542578e+11
         emal production waste
                                 1.859462e+11
         ema2_production_waste
                                 1.587683e+11
         ema3_production_waste
                                 1.465370e+11
         total_releases
                                 1.909209e+10
         on-site_release_total
                                 1.894071e+10
         dtype: float64
         Lowest Variance Cols:
         surface impndmnt
                                 0.0
         ema3 surface impndmnt
                                 0.0
         ema3 underground
                                 0.0
         releases
                                 0.0
         ema3_m71
                                 0.0
         m71
                                 0.0
         m63
                                 0.0
         ema1_releases
                                 0.0
         ema1_underground
                                 0.0
         ema2 releases
                                 0.0
         dtype: float64
         Top 20 Correlations to Age-Adjusted Rate:
         age-adjusted rate
                                  1.000000
         emal_age-adjusted_rate
                                  0.998972
         ema2_age-adjusted_rate
                                  0.998260
         ema3_age-adjusted_rate
                                  0.997837
         ema3_population
                                  0.496632
         ema2_population
                                  0.496499
         emal population
                                  0.496255
         population
                                  0.495891
         ema3 defining PM2.5
                                  0.443200
         ema2 defining PM2.5
                                  0.441238
         ema1_defining_PM2.5
                                  0.436570
         defining_PM2.5
                                  0.422032
                                  0.419695
         aqi
         emal_aqi
                                  0.417988
         ema2 aqi
                                  0.416747
         ema3 aqi
                                  0.416067
         defining Ozone
                                  0.390418
         emal defining_Ozone
                                 0.385056
         ema3 defining NO2
                                 0.384346
         Name: and addited rate dtune: float61
```

All Merged Data w. EMA, Colored by Age-Adjusted Cancer Rates



Algorithm: Predictive Modelling

Now lets see about predicting our values in Y, **life exp, and cancer rates** using our design matrix, **TRI and AQI values**. Since we have multiple such Y columns, we will do so in a **One-vs-AII** approach.

Model Selection

I originally tried a linear model, **ElasticNet** but found that the approach gave very bad R2 values. This is likely because our feature space isn't linear, as we saw in the above scatterplots.

Moreover, we are now considering nearly **500 columns**, many of which will be completely irrelevant for our predictions. I originally tried to use **Backward Featrure Selection** by predicting the <code>age-adjusted_rate</code> values, but this proved computationally intractable since our feature space is so big, as we would have to do this **for every single Y column** when using One-vs-All.

Thus, I decided to use a model with **selects features implicitly**, **Gradient Boosting Regression**. It was the perfect choice since it's a tree method which allows for non-linear relationships, automatically uses only the most relevant features, and returns to me the most important features for each and every Y column I'm trying to predict.

Splitting the Data (Originally src/etl/preprocess.py)

The data will need to split in a stratifed way to prevent poisioning the test set. That is, our DataFrame is **not unique by latitiude and longitude** since we are consdering multiple years for each lat/lon pair. Also, our values are **not on the same scale**, so we will need to scale the values to a normalized range.

The function below will do all of the above, and statify by latitude and longitude for us. It will also reform the index of the DataFrame to be lat, lon, year.

```
In [26]: import json

import geocoder
from sklearn.model_selection import train_test_split
from sklearn preprocessing import MaxAbsScaler
```

```
In [27]: def split(df, y_cols=None, pct=0.20, shuffle=True, normalize=True):
             Splits DataFrame into training and test sets. If `y_cols`
             is provided, also breaks out those columns as a separate
             Numpy Array. Will return either:
                 х, у
                 or
                 x_train, y_train, x_test, y_test
             The split is stratified by Lat and Lon.
             :param df:
                                  DataFrame to split
             :param y_cols:
                                 Y columns of DataFrame
             :param pct:
                                  Validation Percent
             :param shuffle: Shuffle the rows before splitting
             :param normalize
                                  Scale the data using MaxAbsScaler
             :return:
                                  Numpy Arrays
             remove_lat_lon_afterwords = False
             if 'lat' not in df or 'lon' not in df:
                 df['lat'] = [d[0] for d in df.index]
                 df['lon'] = [d[1] for d in df.index]
                 remove_lat_lon_afterwords = True
             df['strat_col'] = df['lat'].astype(str) + " " + df['lon'].astype(str)
             if y cols:
                 y = df[y_cols]
                 cols = [c for c in df.columns if c not in y cols]
                 x = df[cols]
                 x_train, x_test, y_train, y_test = train_test_split(x, y, test_size=pct, s
                                                                       stratify=df['strat col
                 x train = x train.drop('strat col', 1)
                 x_test = x_test.drop('strat_col', 1)
                 if remove lat lon afterwords:
                     x_train = x_train.drop(['lat', 'lon'], 1)
x_test = x_test.drop(['lat', 'lon'], 1)
                 # Fit the scaler to training only
                 if normalize:
                     x_scaler = MaxAbsScaler()
                     x_scaler.fit(x_train)
                     y_scaler = MaxAbsScaler()
                     y_scaler.fit(y_train)
                      return x_scaler.transform(x_train), y_scaler.transform(y_train), \
                             x_scaler.transform(x_test), y_scaler.transform(y_test)
                 else:
                      return x_train, y_train, x_test, y_test
                 # No Y columns provided, simply split the data
                 x_train, x_test = train_test_split(df, test_size=pct, shuffle=shuffle,
                                                      stratify=df['strat col'])
                 x train = x train.drop('strat col', 1)
                 x_test = x_test.drop('strat_col', 1)
                 if remove lat lon afterwords:
                     x_train = x_train.drop(['lat', 'lon'], 1)
                      x_{test} = x_{test.drop(['lat', 'lon'], 1)}
```

Fitting the Model

We will define the function below to fit a One-vs-All Classifier. We will use this function in a loop to fit for every column in Y (Ignoreing EMA columns in Y)

```
In [28]: | from sklearn.ensemble import GradientBoostingRegressor
         from sklearn metrics import mean squared error r? score
In [29]: def fit_predict(x_tr, y_tr, x_te, y_te, label):
              Fits a prediction model in a One-vs-All fashion. Prints
              the results to the console, and returns the classifier object.
              :param x_tr:    X train
:param y_tr:    Y train (1D)
:param x_te:    X test
:param y_te:    Y test (1D)
              :param x tr:
                              X train
              :param label: Label for Y column value
              :return:
                               GBM regressor instance
              gbm = GradientBoostingRegressor(n_estimators=200)
              gbm.fit(x_tr, y_tr)
              print('---', label, '-----')
              for i in sorted(zip(gbm.feature_importances_, x_cols))[-5:]:
                  print(i)
              print("MSE", mean squared error(y te, gbm.predict(x te)))
              print("R2 ", r2_score(y_te, gbm.predict(x_te)))
              return gbm
```

Results

Below are the classification results for every column in Y using a Vanilla GBM with 200 estimators.

```
In [30]: |# Split train and test
         x_train, y_train, x_test, y_test = split(df, y_cols=y_cols)
         # Loop over all Y, ignore EMA columns
         for i in range(len(y_cols)):
             if 'ema' in y_cols[i] or 'sma' in y_cols[i]:
                 continue
            fit predict(x train v train[ il x test v test[ il v cols[il)
         --- population ------
         (0.04395475545428217, 'sector_Miscellaneous Manufacturing')
         (0.049494817251025405, 'm90')
         (0.12245244946747659, 'total_releases')
         (0.20750437389110535, 'sector_Petroleum Bulk Terminals')
         (0.27811621944126635, 'aqi')
         MSE 0.0016256221802137072
         R2 0.9342014901498537
         --- age-adjusted rate ------
         (0.056544920791114395, 'sector_Petroleum Bulk Terminals')
         (0.06701080978716005, 'sector_Tobacco')
(0.08296701667613972, 'lat')
         (0.13865742380446494, 'lon')
         (0.22376649378566038, 'aqi')
         MSE 0.001225445720494672
         R2 0.9302064440473402
         --- cancer_Brain and Other Nervous System ------
         (0.03642829129388199, 'sector_Miscellaneous Manufacturing')
         (0.07387939624873051, 'defining_PM2.5')
```

Explanation of Results, Interpretation

The results are overall quite good! The R2 value is consistently high, but I think it would be interesting to introduce some **value jitter to the X training set** in order to see how robust our findings really are. We can't have a model that only works in delicate circumstances.

Tree Visualization for first three levels of Age-Adjusted Rate GBM (first tree)

```
import io
import pydotplus
from PIL import Image
from matplotlib import pyplot as plt
from sklearn.externals.six import StringIO
from sklearn.tree import DecisionTreeRegressor
from sklearn tree import export graphyiz
/home/cole/.pyvenv/lib/python3.7/site-packages/sklearn/externals/six.py:31: Depr
ecationWarning: The module is deprecated in version 0.21 and will be removed in
version 0.23 since we've dropped support for Python 2.7. Please rely on the offi
cial version of six (https://pypi.org/project/six/).
    "(https://pypi.org/project/six/).", DeprecationWarning)
```

```
In [32]: def visualize_best_tree(x, y, levels=3):
                t = GradientBoostingRegressor(max_depth=levels, n_estimators=200)
               t.fit(x, y)
               t = t.estimators_{0}, 0
               dot_data = StringIO()
               export_graphviz(t, out_file=dot_data,
                                   filled=True, rounded=True,
                                   special characters=True)
               graph = pydotplus.graph_from_dot_data(dot_data.getvalue())
               img = Image.open(io.BytesIO(graph.create png()))
                img.save('doc/plots/age-adjusted rate-tree-viz.png')
                disnlav(ima)
In [33]: visualize hest tree(x train v train[ · 11)
In [34]: print("Legend:")
          print("X8:", x_cols[8])
print("X2:", x_cols[2])
print("X66:", x_cols[66])
print("X40:", x_cols[40])
print("X1:", x_cols[41])
          print("X1:", x_cols[1])
print("X70:", x_cols[70])
print("X61:" x_cols[61])
           Legend:
          X8: defining_PM2.5
          X2: lon
          X66: off-site_contain
          X40: m65
          X1: lat
           X70: sector_Chemical Wholesalers
           X61: off-site_treated_total
          Adding Noise to Training Data
In [35]: imnort numny as no
```

```
In [36]: noise = np.random.normal(0,1, x_train.shape) * 0.01
         x train += nnise
```

```
In [37]: # Loop over all Y, ignore EMA columns
         for i in range(len(y_cols)):
             if 'ema' in y_cols[i] or 'sma' in y_cols[i]:
                 continue
             fit nredict(x train v train(: il. x test v test(: il. v cols[il)
         --- population ------
         (0.042516741454621206, 'defining_NO2')
         (0.04552335453346455, 'm90')
         (0.07591751218152834, 'sector_Computers and Electronic Products')
         (0.20502468313661323, 'sector_Petroleum Bulk Terminals')
         (0.27127832321126116, 'agi')
         MSE 0.0017296458852134795
         R2 0.9299910377695986
         --- age-adjusted rate ------
         (0.06482359165831117, 'sector_Tobacco')
         (0.06546907031610365, 'sector_Petroleum Bulk Terminals')
         (0.09541190645495926, 'lon')
         (0.09560436925002326, 'lat')
         (0.23751731058954678, 'agi')
         MSE 0.0019030452661857767
         R2 0.891614704719548
         --- cancer_Brain and Other Nervous System ------
         (0.04844408768102432, 'sector_Electrical Equipment')
         (0.06695685036118884, 'sector Computers and Electronic Products')
```

Evaluation: Goodness of Analytic

We can see that adding noise may not help our prediction, but it doesn't destroy it either. This leads me to believe that the predictions we originally have are reasonably accurate at least on the surface.

If the model were simply fitting to, and memorizing the input data, then the noise would have destoryed the predictive accuracy (R2 value) of our model more than it did.

Also in-line with our earlier correlation analysis, no one feature is important alone in predicting cancer rates and life expectancy. They all contribute a fair amount, without overfitting to one given feature. We see that in our GBM feature_importances values, no one column really exceeds about 0.35 for any analysis. (This value ranges from 0.0 to 1.0, and all values add up to 1.0)

Evaluation: Findings

More often than not, the most defining characteristics for different cancer rates were the number of **Petroleum Bulk Terminals** dumpings, and the **Air Quality Index** for a given area.

Surprisingly, the defining characteristic for the AQI that was most important in determing cancer rates was **PM2.5**. From previous experience in Air Quality research, PM2.5 is often **not** the defining characteristic. Normally it is Ozone, which has been clinically proven to have an adverse affect on sensitive groups.

I also intentionally left **lat** and **lon** as features of the data set to see if cancer rates were simply concentrated to parts of the United States, rather than factors *within* those areas. In this way, these features model the **spurious correlation** of our other features, since latitude and longitude were part of our grouping index.

In age-adjusted_rate we see that lat and lon are both in the top 5 most important features for the GBM regressor which probably indicates some of these relationships are spurious.

Evaluation: Future Work

Future work would need to be conducted in determining the causal relationships between these features and actual cancer rates. It is not enough to say that these correlations are linked. However, these studies are often expensive and the novelty of this application is to show which factors are worth exploring. For that, I believe Air Quality and Tobacco / Petroleum Dumping should be explored in greater detail to assess the health risk as it relates to life expectancy and cancers.

Moreover, more work can be done to this application to better relate the features in the design matrix. For instance, not all types of dumping may be equally harmful, and this model does not currently have a way to determine that as it sees the *kind of* dumping as an independent feature to the *industry* that is dumping. This however requires an intuitive understanding and industry knowledge of toxic dumping.

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In I		