Visualizing high dimensional TEDS data

In [2]:
```
import matplotlib.pyplot as plt
import pandas as pd

import seaborn as sns

import numpy as np
import scipy.stats as ss

from sklearn.decomposition import PCA
from sklearn.preprocessing import StandardScaler
```

In [7]:
```
df=pd.read_csv('__TEDSCountyLevel2008-2017_2.csv')
```

In [3]:
```
%matplotlib inline
```

In [8]:
```
df.head(10)
```

Out[8]:

	Alcohol	Amphetamine	Barbiturates	Benzine	Age	State	Cocaine	County Name	Gender	Hal	...	Pr
0	0	0	0	0	18-24	Indiana	0	Pulaski	Male	0	...	0
1	1	0	0	0	55+	Indiana	0	Pulaski	Male	0	...	0
2	1	0	0	0	18-24	Indiana	0	Pulaski	Male	0	...	0
3	0	0	0	0	55+	Indiana	0	Pulaski	Male	0	...	0
4	0	0	0	0	55+	Indiana	0	Pulaski	Male	0	...	0
5	1	0	0	0	25-34	Indiana	0	Pulaski	Male	0	...	0
6	1	0	0	0	35-44	Indiana	0	Pulaski	Male	0	...	0
7	1	0	0	0	25-34	Indiana	0	Pulaski	Male	0	...	0
8	1	0	0	0	0-18	Indiana	0	Pulaski	Male	0	...	0
9	1	0	0	0	25-34	Indiana	0	Pulaski	Male	0	...	0

10 rows × 50 columns

In [9]:
```
#Limit the number of features
features = ['Year','County Name','Gender','Race','Age','Alcohol','Heroin','Marijuana','O
pioid','Cocaine','Other Drugs','Pain Killers', 'Methamphetamine', 'Synthetic Drugs']
df_subset = df[features]
df_subset.head()
```

Out[9]:

	Year	County Name	Gender	Race	Age	Alcohol	Heroin	Marijuana	Opioid	Cocaine	Other Drugs	Pain Killers	M
0	2009	Pulaski	Male	White	18-24	0	0	1	0	0	0	0	0
1	2009	Pulaski	Male	White	55+	1	0	0	0	0	0	0	0
2	2009	Pulaski	Male	White	18-24	1	0	1	0	0	0	0	0
3	2009	Pulaski	Male	White	55+	0	0	1	0	0	0	0	0
4	2009	Pulaski	Male	White	55+	0	0	1	0	0	0	0	0

In [10]:
```
import altair as alt
alt.renderers.enable('notebook')
```

Out[10]:
```
RendererRegistry.enable('notebook')
```

In [11]:
```
#Create a dataset for the 9 substances to use in correlations
features = ['Alcohol','Heroin','Marijuana','Opioid','Cocaine','Other Drugs','Pain Killer
s', 'Methamphetamine', 'Synthetic Drugs']
df_only_features = df_subset[features]
df_only_features.head()
```

Out[11]:

	Alcohol	Heroin	Marijuana	Opioid	Cocaine	Other Drugs	Pain Killers	Methamphetamine	Synthetic Drugs
0	0	0	1	0	0	0	0	0	0
1	1	0	0	0	0	0	0	0	0
2	1	0	1	0	0	0	0	0	0
3	0	0	1	0	0	0	0	0	0
4	0	0	1	0	0	0	0	0	0

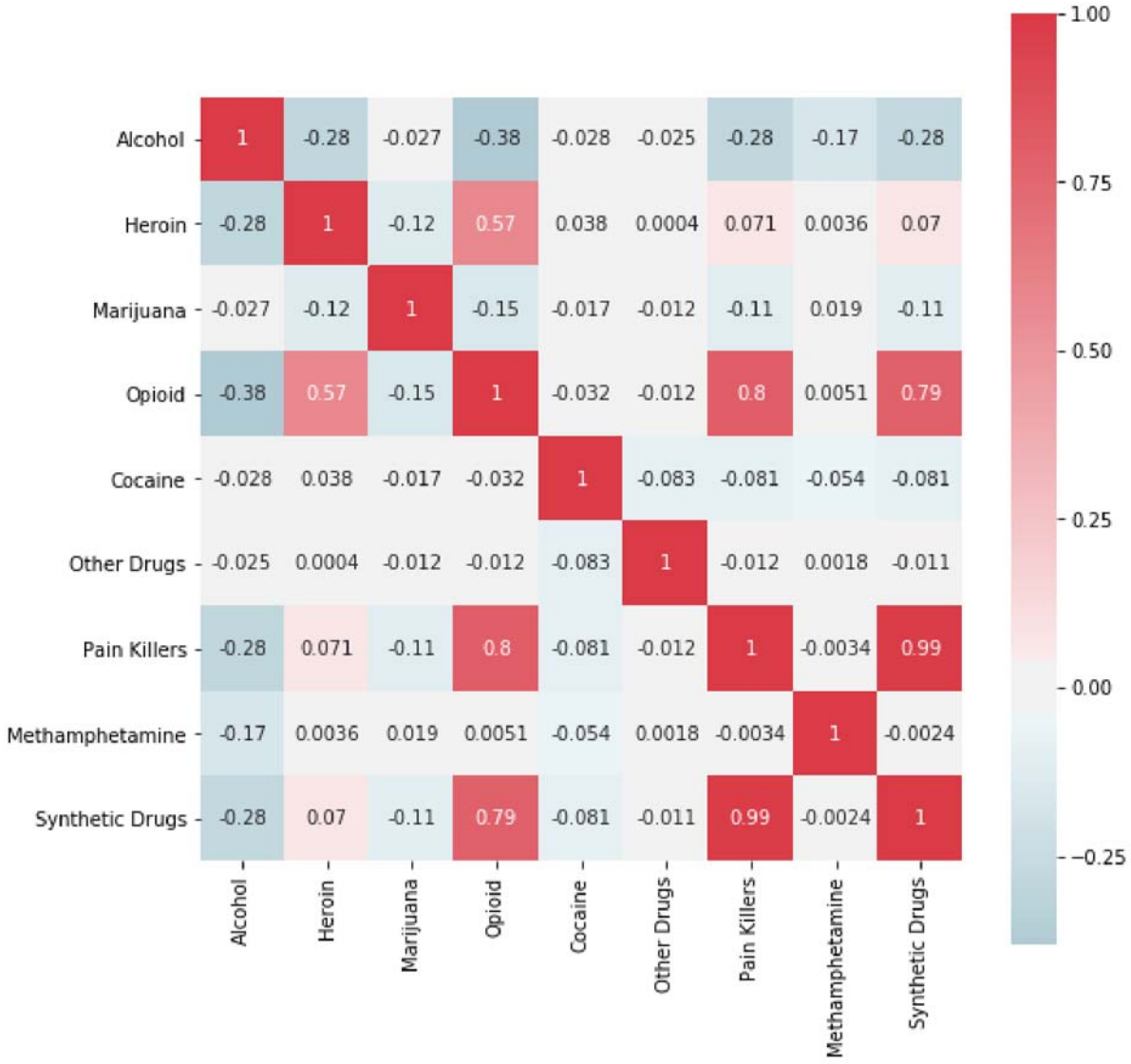
In [12]:
```
n_samples, n_features = df_only_features.shape
print(n_samples)
print(n_features)
```

```
335866
9
```

In [11]:
```
#Check correlation between substances
import seaborn as sns

f, ax = plt.subplots(figsize=(9, 9))
corr = df_only_features.corr()
sns.heatmap(corr, mask=np.zeros_like(corr, dtype=np.bool), cmap=sns.diverging_palette(220, 10, as_cmap=True), center=0,
            square=True, ax=ax,annot=True)
```

Out[11]: <matplotlib.axes._subplots.AxesSubplot at 0x1bd9f9260b8>



There is strong correlation between Pain Killers and Synthetic Drugs, supported by the Opioids. Heroin appears in 57% of the cases with Opioids. Alcohol is slightly negatively correlates with all four of them.

Remarkably, there appears to be no appreciable correlation between opioid use and such commonly accepted and complementary socio-economic markers as alcohol, marijuana, or methamphetamine use. In fact, these three factors are likely to be biased towards older population (alcohol), younger population (marijuana) and underprivileged rural population (methamphetamine).

Thus, it appears that opioid abuse have roots that are different from the common socio-economic maladies reflected in these three markers.

In [13]:
```
df_subset['County Name'].nunique()
```

Out[13]:
```
93
```

In [13]:
```
df_subset['Year'].nunique()
```

Out[13]:
```
10
```

In [14]:
```
# Total number of records
len(df_subset)
```

Out[14]:
```
335866
```

In [15]:
```
#How many records per year
df_subset['Year'].value_counts()
```

Out[15]:
```
2017    37459
2015    34739
2011    34677
2014    34470
2012    33984
2013    33616
2016    33170
2009    32049
2010    31827
2008    29875
Name: Year, dtype: int64
```

In [15]:
```
#How many records per county
df_subset['County Name'].value_counts()
```

Out[15]:

Marion	41678
Lake	25329
Allen	16940
Vanderburgh	14085
Monroe	13842
Saint Joseph	13359
Delaware	10453
Madison	9594
Elkhart	8309
Hamilton	8033
Vigo	7895
Howard	6235
Bartholomew	6037
Porter	5455
LaPorte	5086
Morgan	5057
Lawrence	4557
Tippecanoe	4514
Wayne	4391
Grant	4241
Dearborn	3981
Clark	3857
Hendricks	3488
Knox	3350
Noble	3024
Montgomery	3006
Warrick	2969
Kosciusko	2946
Jefferson	2888
Johnson	2834
...	...
Perry	1407
Rush	1281
Shelby	1246
Vermillion	1241
Wells	1195
Jay	1193
Orange	1174
Franklin	1115
Huntington	1107
Parke	1084
Whitley	1031
Carroll	967
Brown	922
Pulaski	906
Jasper	902
Sullivan	812
Blackford	759
Washington	681
Fountain	662
Martin	638
Tipton	563
Switzerland	562
Crawford	511
Harrison	461
Pike	433
Benton	373
Newton	315
Ohio	293
Union	284

```
Warren            229
Name: County Name, Length: 93, dtype: int64
```

In [16]:
```
# General method to variably group the data
def group_dataset(df, col1, col2):
    df_grouped=df.groupby([col1,col2])
    return df_grouped
```

In [17]:
```
#Group by Year and County
df_grouped=group_dataset(df_subset, 'Year','County Name')

#Excluded pain killers and synthetic drugs as highly correclated with opioids
feature=['Alcohol','Heroin','Marijuana','Opioid','Cocaine','Other Drugs','Methamphetamin
e']
df_grouped=df_grouped[feature].mean()

df_grouped.head()
```

Out[17]:

		Alcohol	Heroin	Marijuana	Opioid	Cocaine	Other Drugs	Methamphetamine
Year	**County Name**							
2008	**Adams**	0.905263	0.021053	0.684211	0.105263	0.231579	0.0	0.010526
	Allen	0.863495	0.012480	0.641186	0.048362	0.325273	0.0	0.017161
	Bartholomew	0.651106	0.004914	0.542998	0.226044	0.255528	0.0	0.299754
	Benton	0.840000	0.000000	0.520000	0.040000	0.160000	0.0	0.040000
	Blackford	0.833333	0.000000	0.620370	0.157407	0.175926	0.0	0.009259

In [18]:
```
#Group by County
df_grouped_c=group_dataset(df_subset,'County Name', 'Year')

#exclude pain killers and synthetic drugs as highly correclated with opioids
feature=['Alcohol','Heroin','Marijuana','Opioid','Cocaine','Other Drugs','Methamphetamin
e']
df_grouped_c=df_grouped_c[feature].mean()

df_grouped_c.head()
```

Out[18]:

		Alcohol	Heroin	Marijuana	Opioid	Cocaine	Other Drugs	Methamphetamine
County Name	**Year**							
Adams	**2008**	0.905263	0.021053	0.684211	0.105263	0.231579	0.000000	0.010526
	2009	0.877863	0.007634	0.549618	0.099237	0.175573	0.000000	0.030534
	2010	0.613497	0.141104	0.496933	0.300613	0.184049	0.147239	0.042945
	2011	0.655405	0.141892	0.540541	0.331081	0.155405	0.108108	0.060811
	2012	0.756098	0.073171	0.528455	0.203252	0.178862	0.138211	0.040650

In [19]:
```
# Group by Year to genralize the output
df_grouped_year1=df_subset.groupby('Year')
df_grouped_year1=df_grouped_year1[feature].mean()

df_grouped_year1.head()
```

Out[19]:

	Alcohol	Heroin	Marijuana	Opioid	Cocaine	Other Drugs	Methamphetamine
Year							
2008	0.661891	0.038226	0.506644	0.155213	0.223397	0.000000	0.087464
2009	0.682798	0.049237	0.514837	0.177821	0.194359	0.000000	0.091235
2010	0.632168	0.055833	0.490998	0.186006	0.157602	0.119553	0.096962
2011	0.606050	0.068662	0.483837	0.220607	0.146899	0.138593	0.103556
2012	0.590837	0.094191	0.471398	0.275012	0.156809	0.154926	0.112612

In [20]:
```
# General method to pivot by two predictors
def create_pivot_df(df, col_name, index_name, val_name):
    pivot_df = df.pivot_table(columns=col_name, index= index_name, values= val_name, agg
func='mean')
    return pivot_df
```

In [21]:
```
#create a pivot table for alcohol
create_pivot_df(df_grouped,col_name='Year',index_name='County Name', val_name='Alcohol')
.head()
```

Out[21]:

Year	2008	2009	2010	2011	2012	2013	2014	2015	2
County Name									
Adams	0.905263	0.877863	0.613497	0.655405	0.756098	0.812030	0.805195	0.697143	0.559
Allen	0.863495	0.834094	0.812057	0.794048	0.672963	0.736813	0.729648	0.674190	0.665
Bartholomew	0.651106	0.605505	0.442708	0.400815	0.458841	0.522551	0.501475	0.410745	0.345
Benton	0.840000	0.850000	0.684211	0.794118	0.852941	0.659091	0.708333	0.750000	0.780
Blackford	0.833333	0.787234	0.216216	0.523810	0.519231	0.386364	0.440678	0.454545	0.378

In [22]:
```
def draw_heatmap(df, nrows, ncols):
    f, ax = plt.subplots(figsize=(nrows, ncols))
    sns.heatmap(df,  cmap=sns.cubehelix_palette(8, start=.5, rot=-.75),ax=ax,xticklabels
=2, square=True, vmin=0, vmax=.9)
```

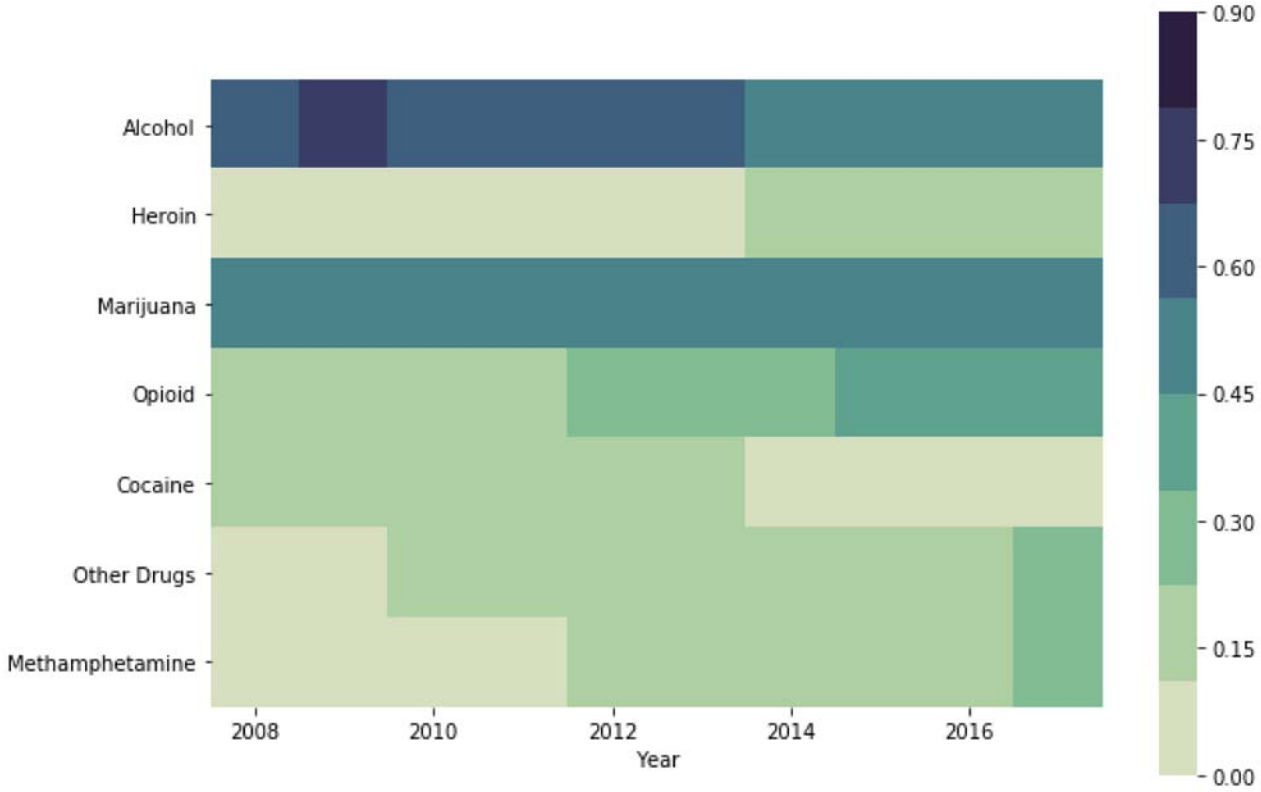
In [23]:
```
def draw_heatmap_corr(df, nrows, ncols):
    f, ax = plt.subplots(figsize=(nrows, ncols))
    df=df.corr()
    sns.heatmap(df,  cmap=sns.diverging_palette(220, 10, as_cmap=True), center=0,ax=ax,
square=True,annot=True,cbar_kws={"shrink": .5})
```

In [24]:
```
#Transpose the matrix
df_grouped_year_T= df_grouped_year1.T
df_grouped_year_T
```

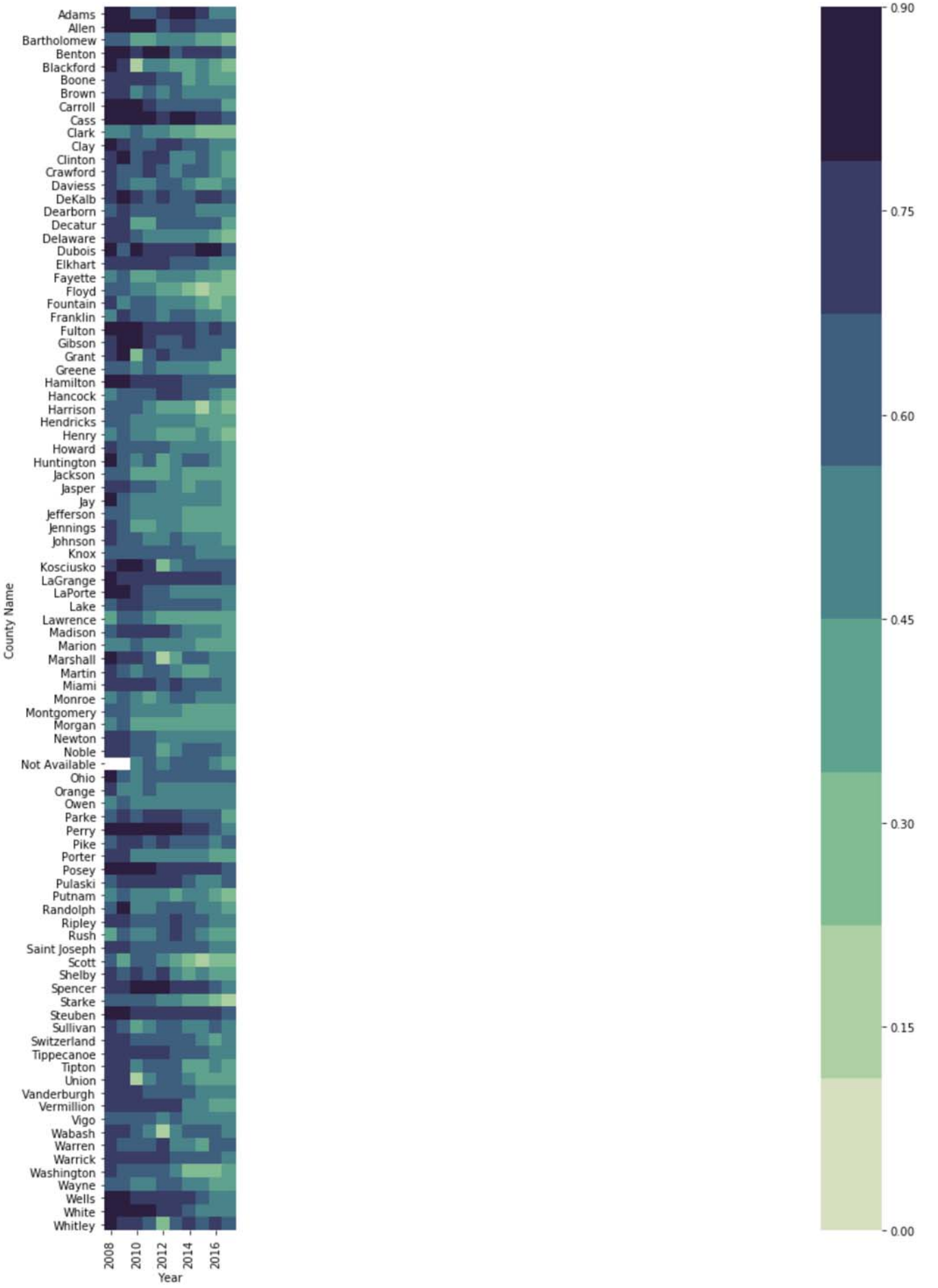
Out[24]:

Year	2008	2009	2010	2011	2012	2013	2014	2015
Alcohol	0.661891	0.682798	0.632168	0.606050	0.590837	0.585792	0.559066	0.528743
Heroin	0.038226	0.049237	0.055833	0.068662	0.094191	0.105396	0.121787	0.156625
Marijuana	0.506644	0.514837	0.490998	0.483837	0.471398	0.476172	0.478648	0.483462
Opioid	0.155213	0.177821	0.186006	0.220607	0.275012	0.303248	0.314244	0.339273
Cocaine	0.223397	0.194359	0.157602	0.146899	0.156809	0.131277	0.112446	0.106825
Other Drugs	0.000000	0.000000	0.119553	0.138593	0.154926	0.165516	0.177894	0.202654
Methamphetamine	0.087464	0.091235	0.096962	0.103556	0.112612	0.121995	0.141514	0.159936

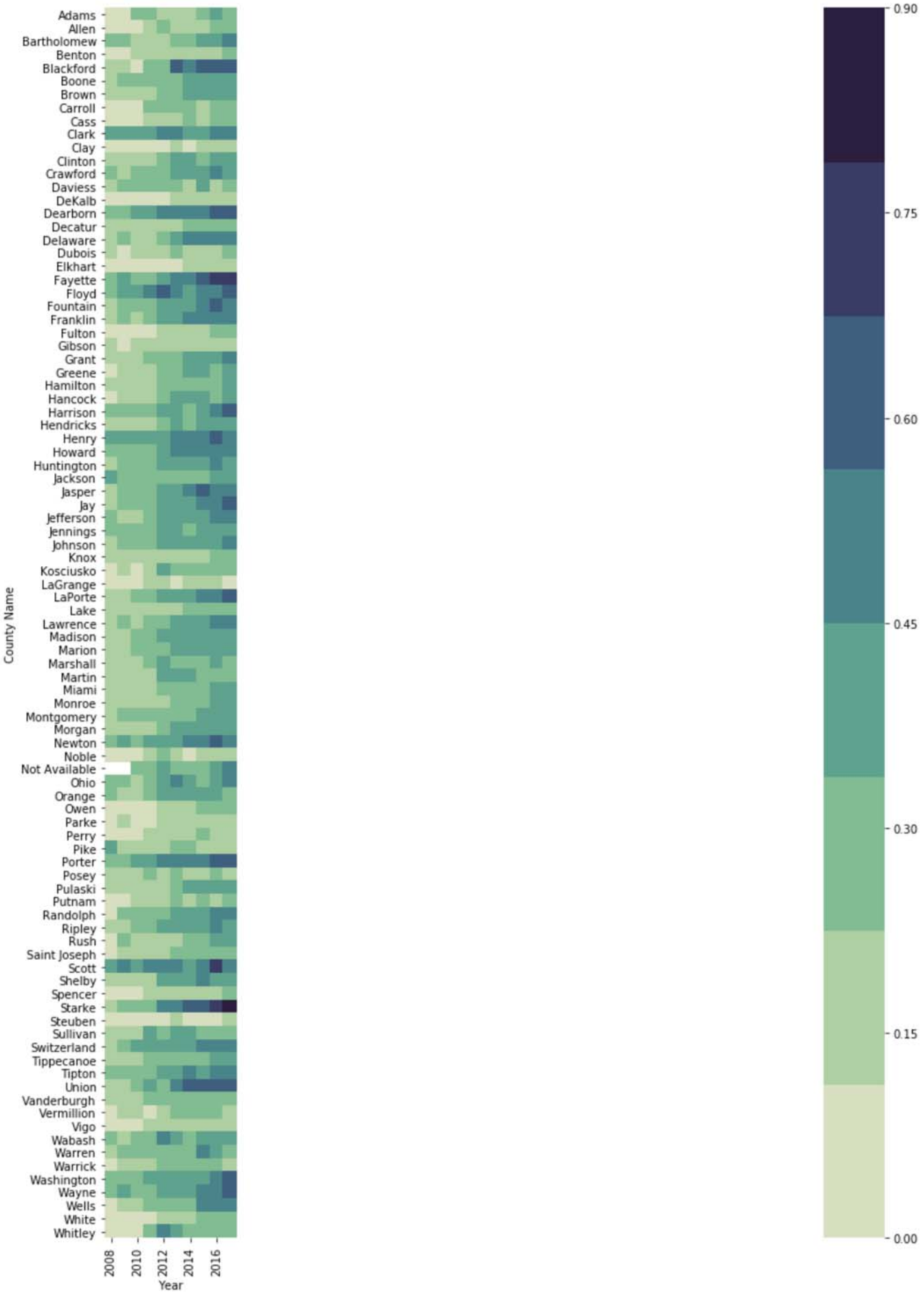
In [27]:
```
# Show a trend relative to other substances
draw_heatmap(df_grouped_year_T, 10, 7)
```



In [28]:
```
# Draw a heatmap by County for Alcohol
draw_heatmap(
create_pivot_df(df_grouped,'Year','County Name', 'Alcohol'), 186, 20)
```

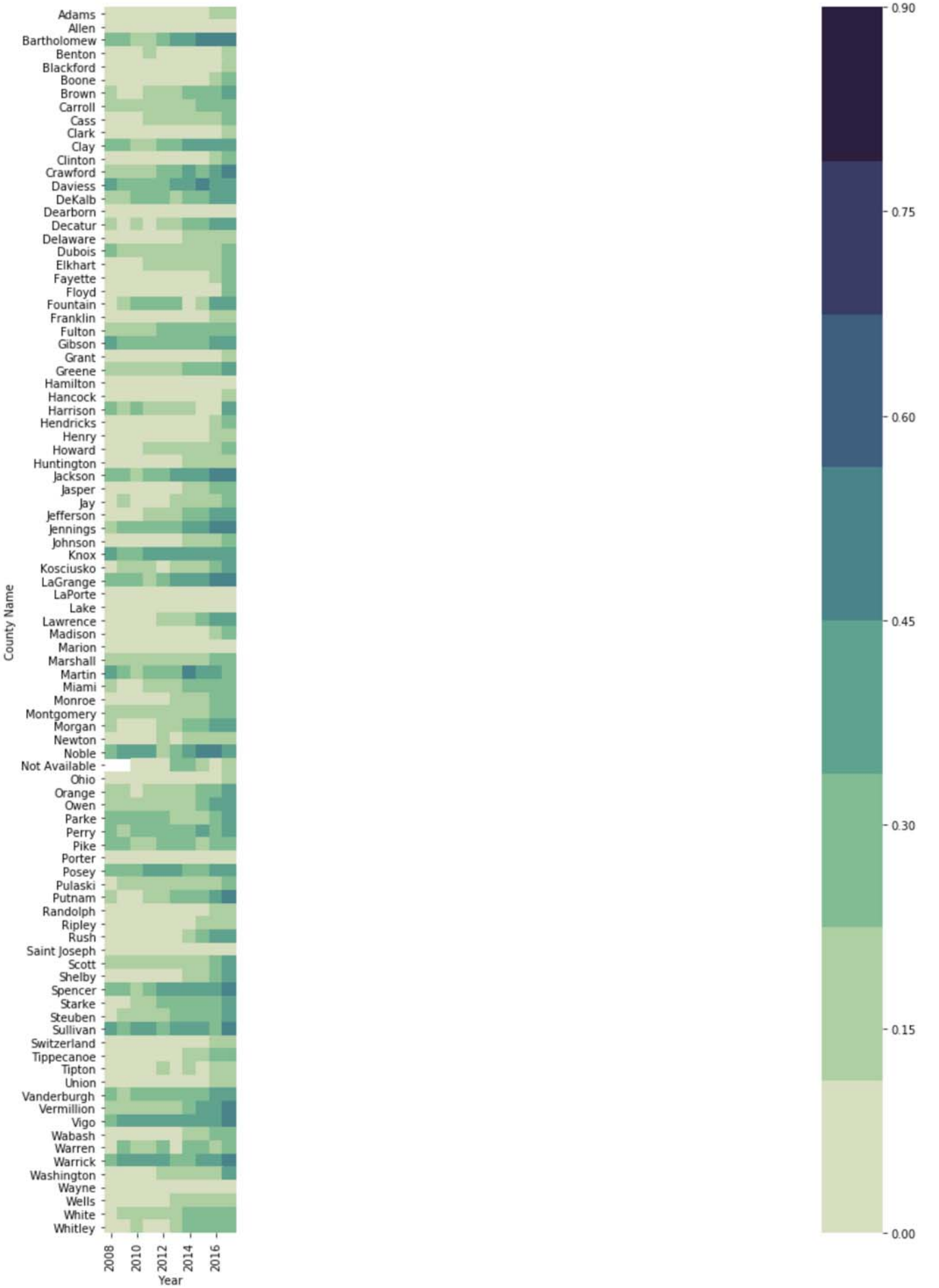


In [29]:
```
#Opioid
draw_heatmap(create_pivot_df(df_grouped,'Year','County Name', 'Opioid'), 186, 20)
```



In [30]:
```
#Marijuana
draw_heatmap(create_pivot_df(df_grouped,'Year','County Name', 'Marijuana'), 186, 20)
```

In [31]:
```
# Methamphetamine
draw_heatmap(create_pivot_df(df_grouped,'Year','County Name', 'Methamphetamine'), 186, 20)
```



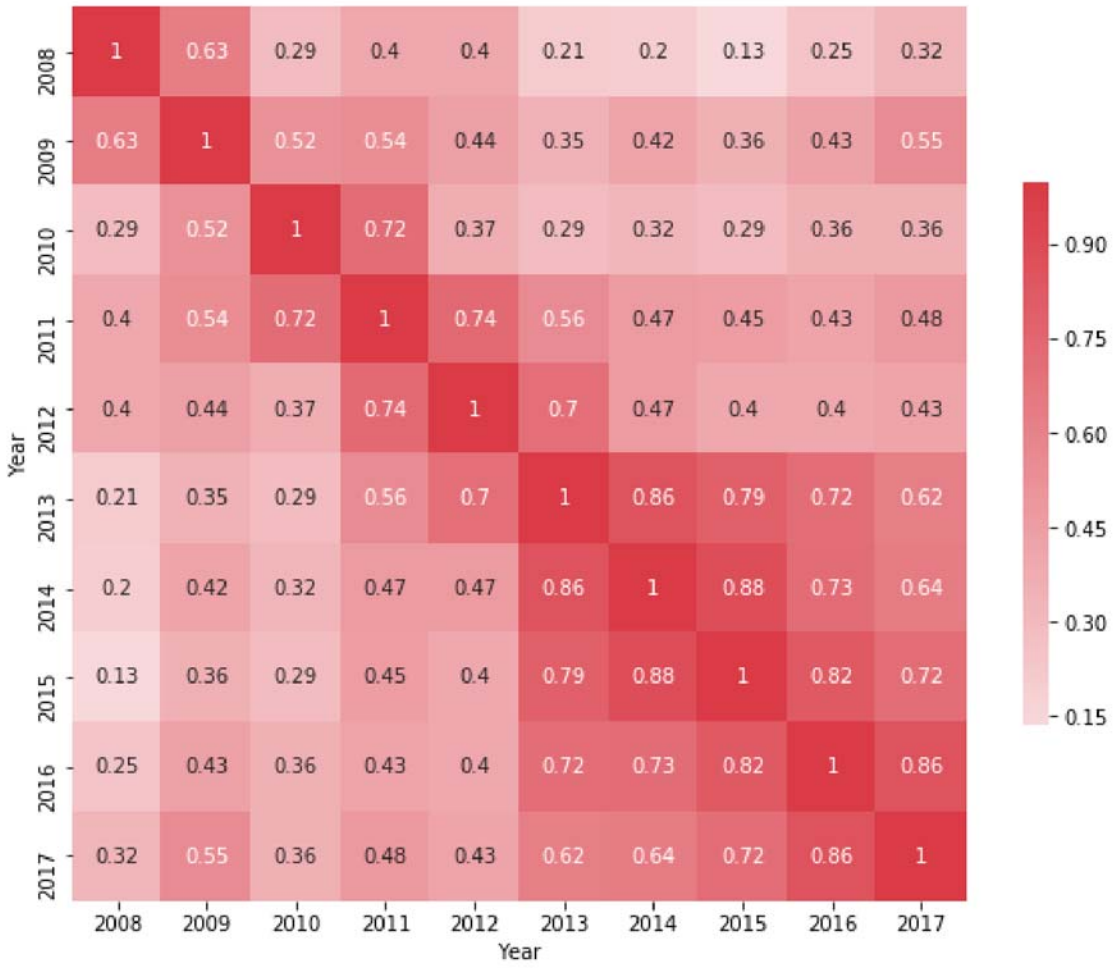
From the four heatmaps above, one could infer the following observations:

- while Alcohol remains the most commonly abused substance, its dominance is vaning as the prevalence of other substances, especially opiods and methamphethamine increases. At the same time, marijuana abuse remains fairly stable.

Suplimentary Analysis: Identify any years that differ from the majority

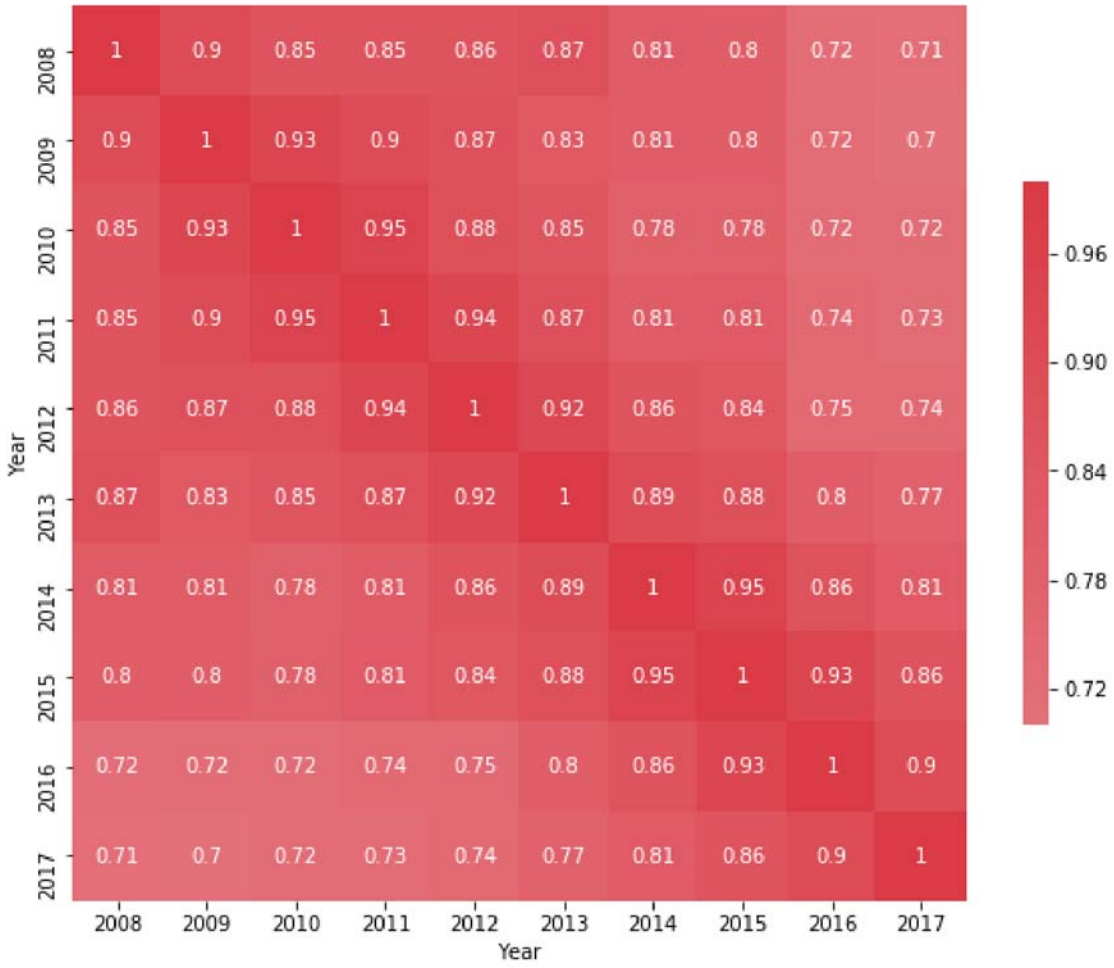
In [ ]:
```
draw_heatmap_corr(create_pivot_df(df_grouped,'Year','County Name', 'Alcohol'), 10,10)
```

In [33]:
```
#Marijuana
draw_heatmap_corr(create_pivot_df(df_grouped,'Year','County Name', 'Marijuana'), 10,10)
```



In [34]:
```
#Opioids
draw_heatmap_corr(create_pivot_df(df_grouped,'Year','County Name', 'Opioid'), 10,10)
```

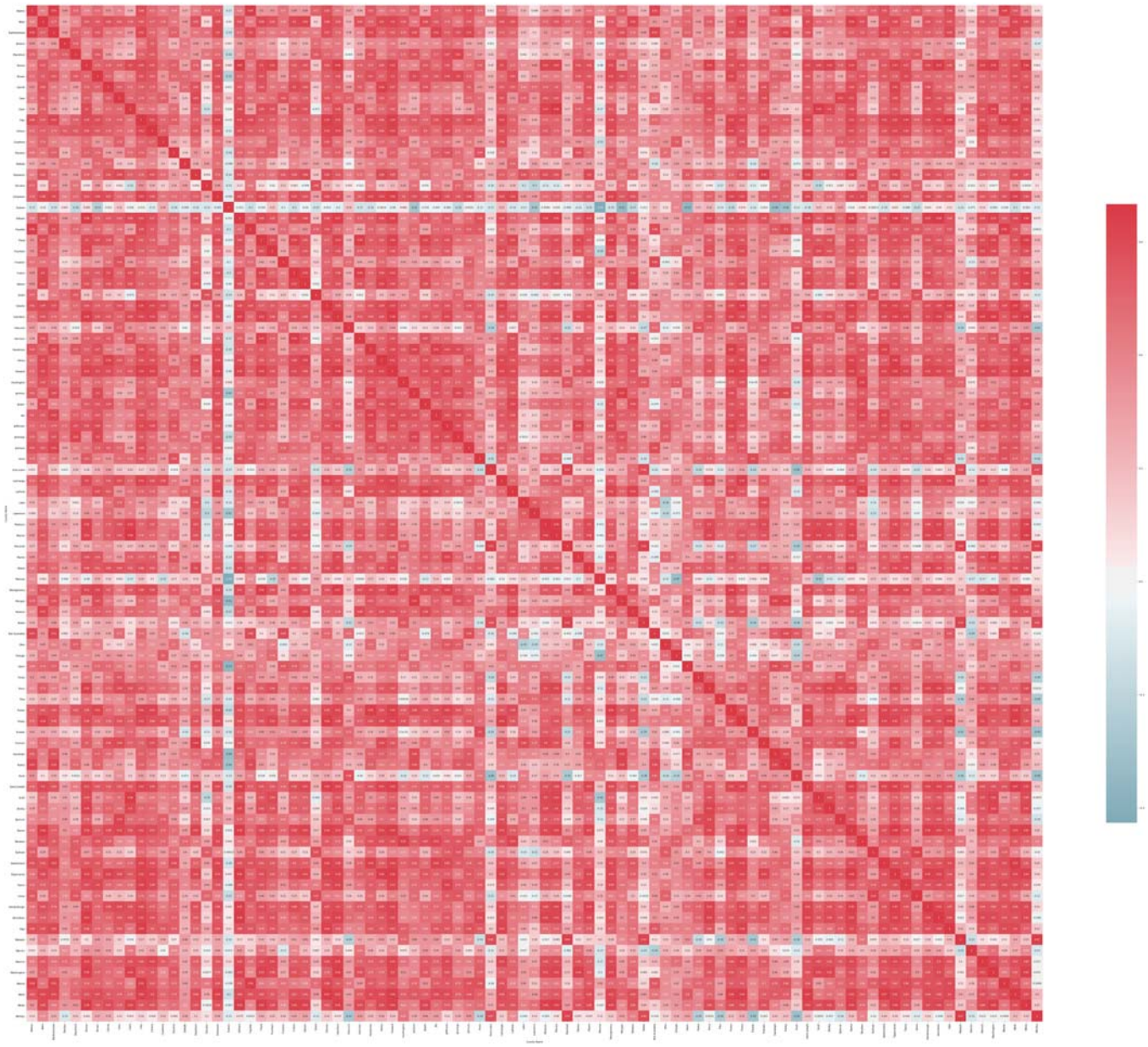
In [35]:
```
#Methamphetamine
draw_heatmap_corr(create_pivot_df(df_grouped,'Year','County Name', 'Methamphetamine'), 10,10)
```



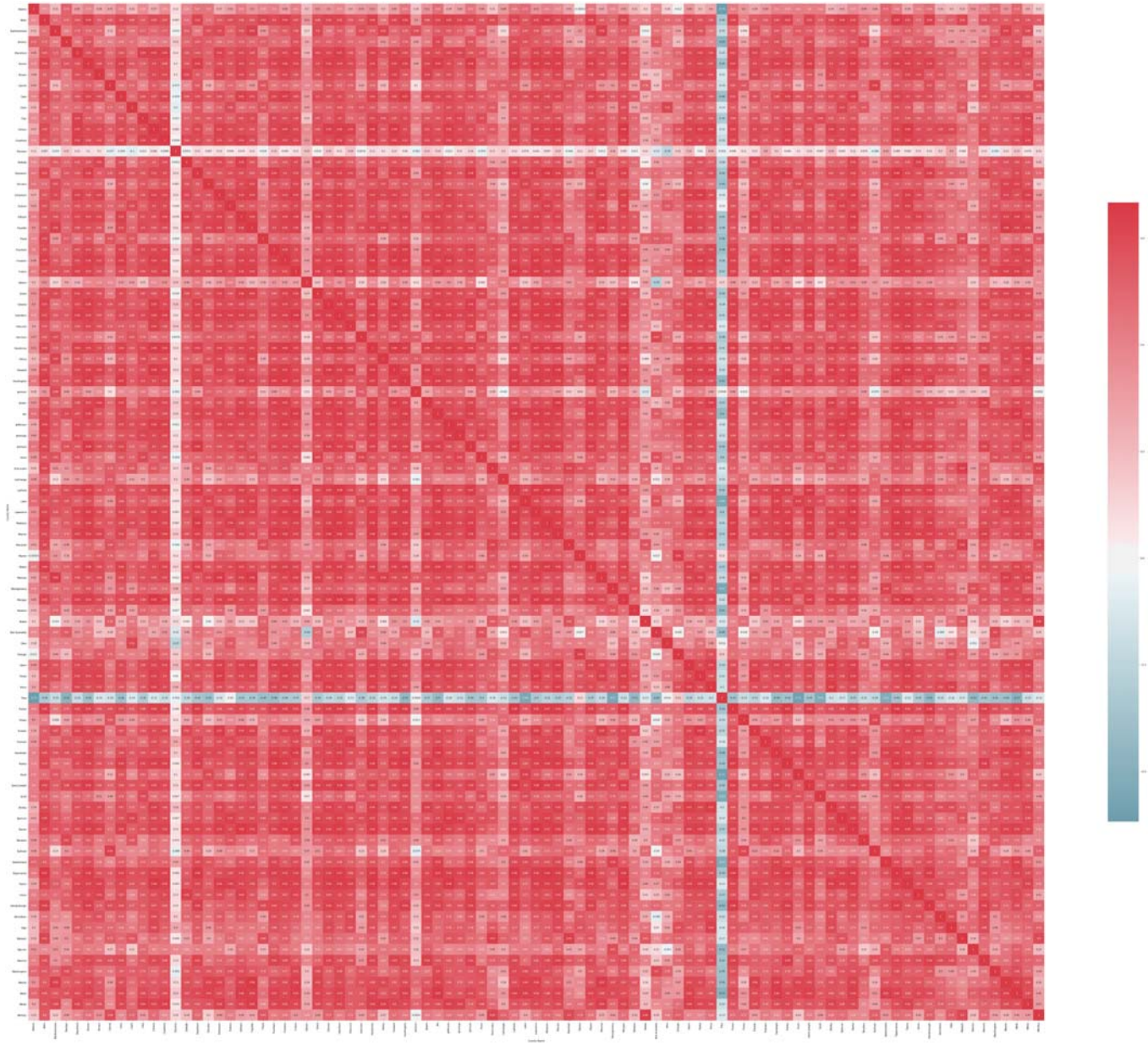
Beginning 2012 the usage patterns became very similar for the malority of the subjects that requested treatment in Indiana.

Suplimentary Analysis: Identify counties that differ from the majority

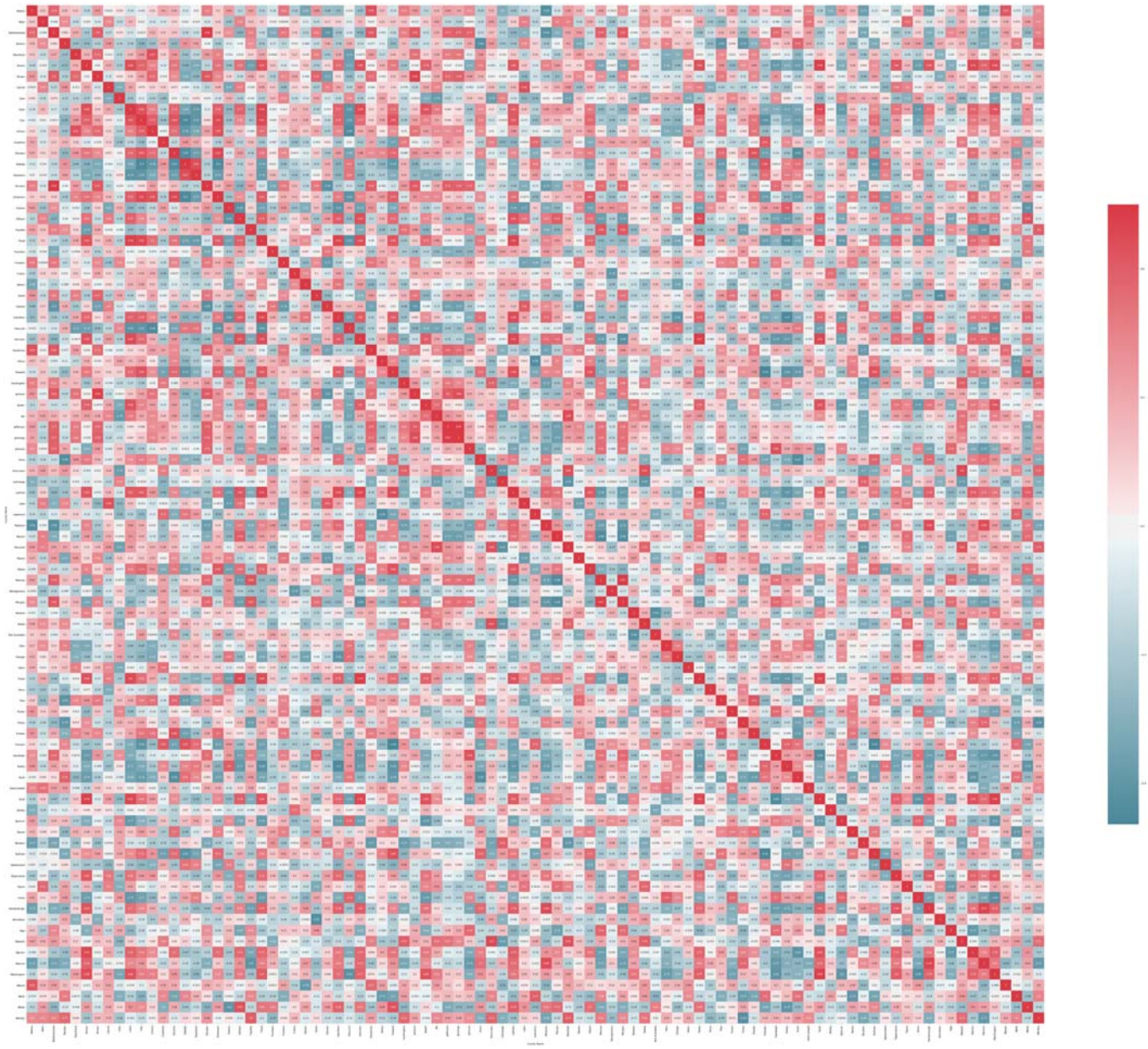
In [37]:
```
draw_heatmap_corr(create_pivot_df(df_grouped,'County Name','Year', 'Alcohol'), 93,93)
```



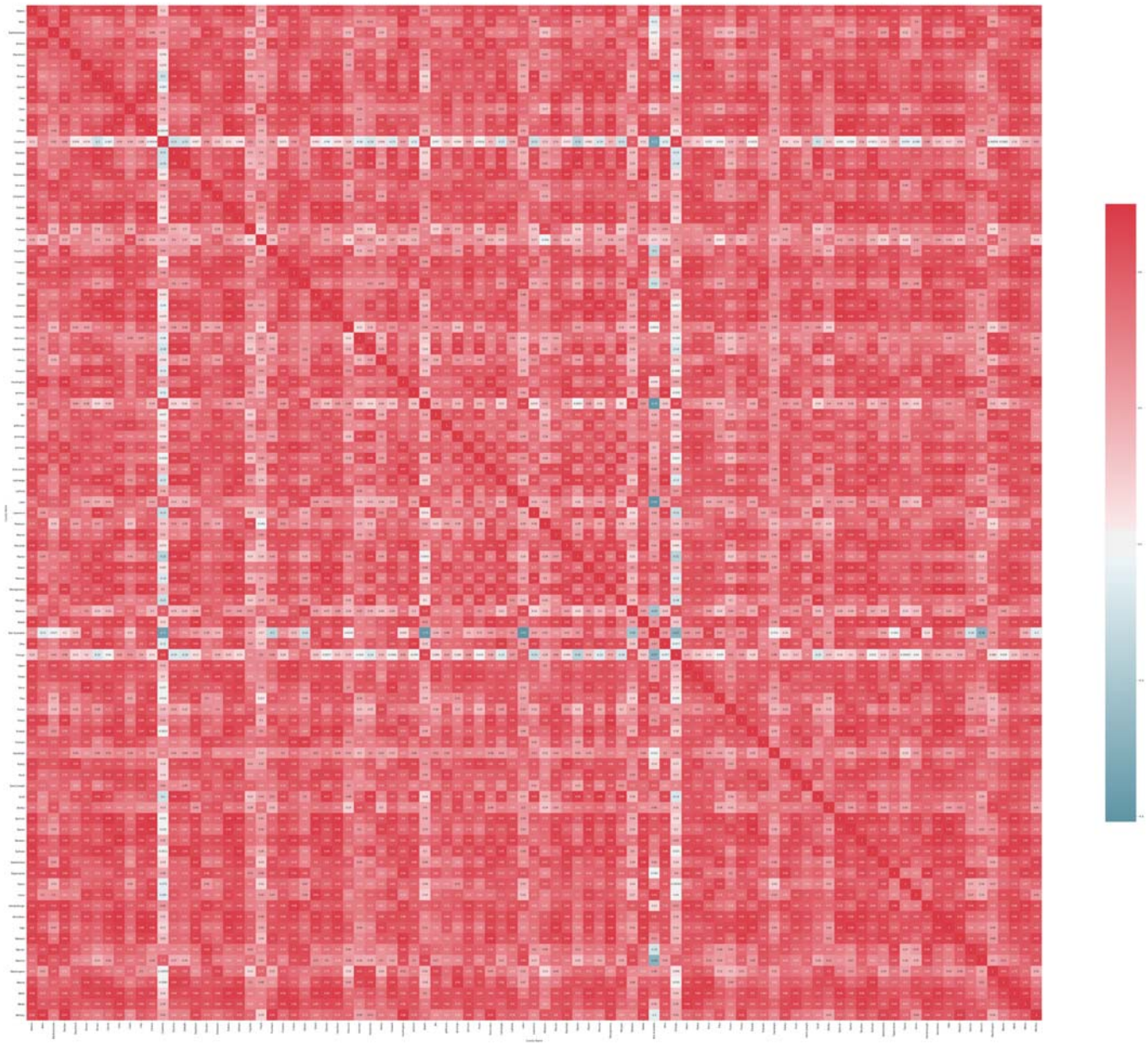
In [38]:
```
draw_heatmap_corr(create_pivot_df(df_grouped,'County Name','Year', 'Opioid'), 93,93)
```



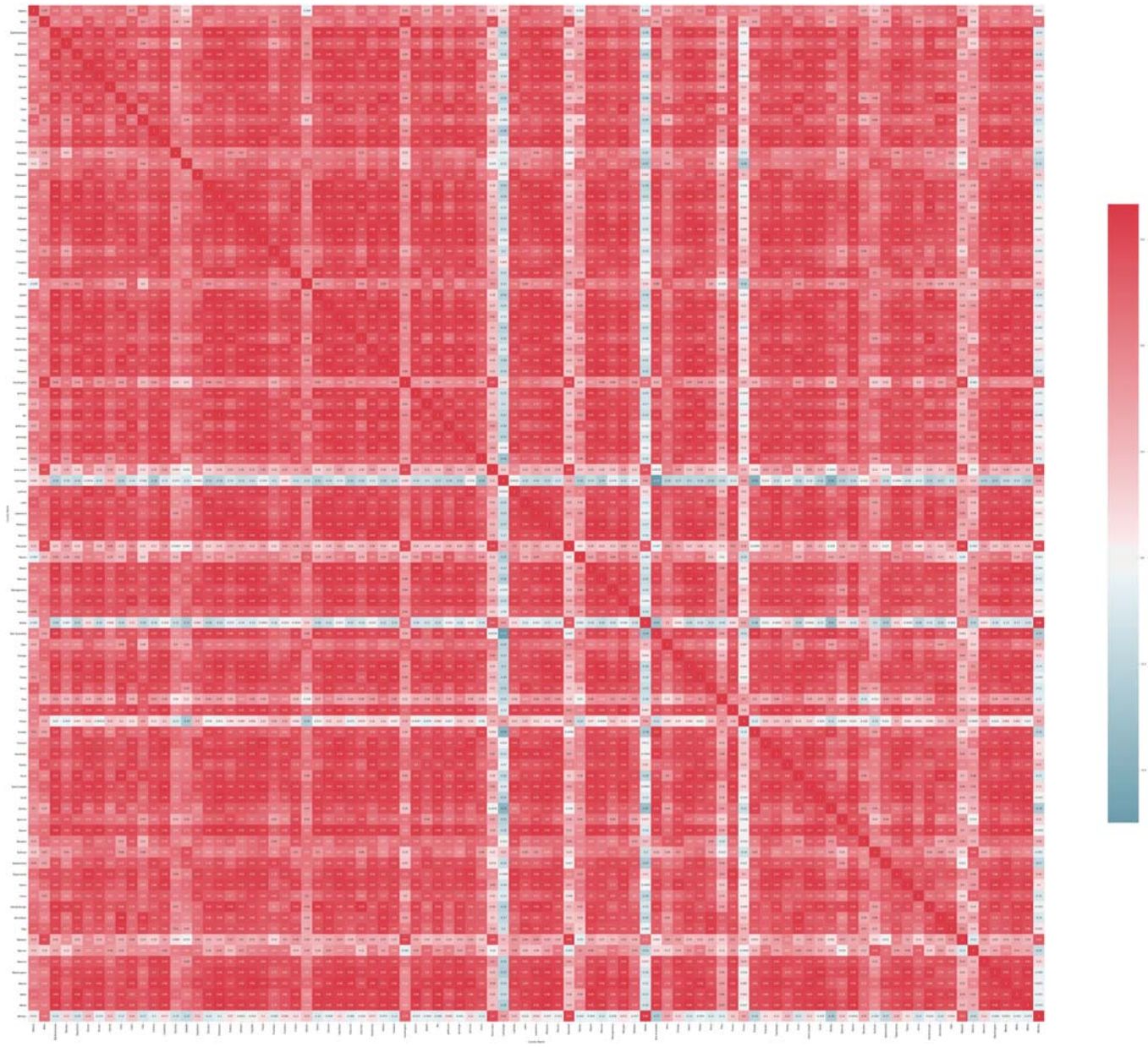
In [39]:
```
draw_heatmap_corr(create_pivot_df(df_grouped,'County Name','Year', 'Marijuana'), 93,93)
```



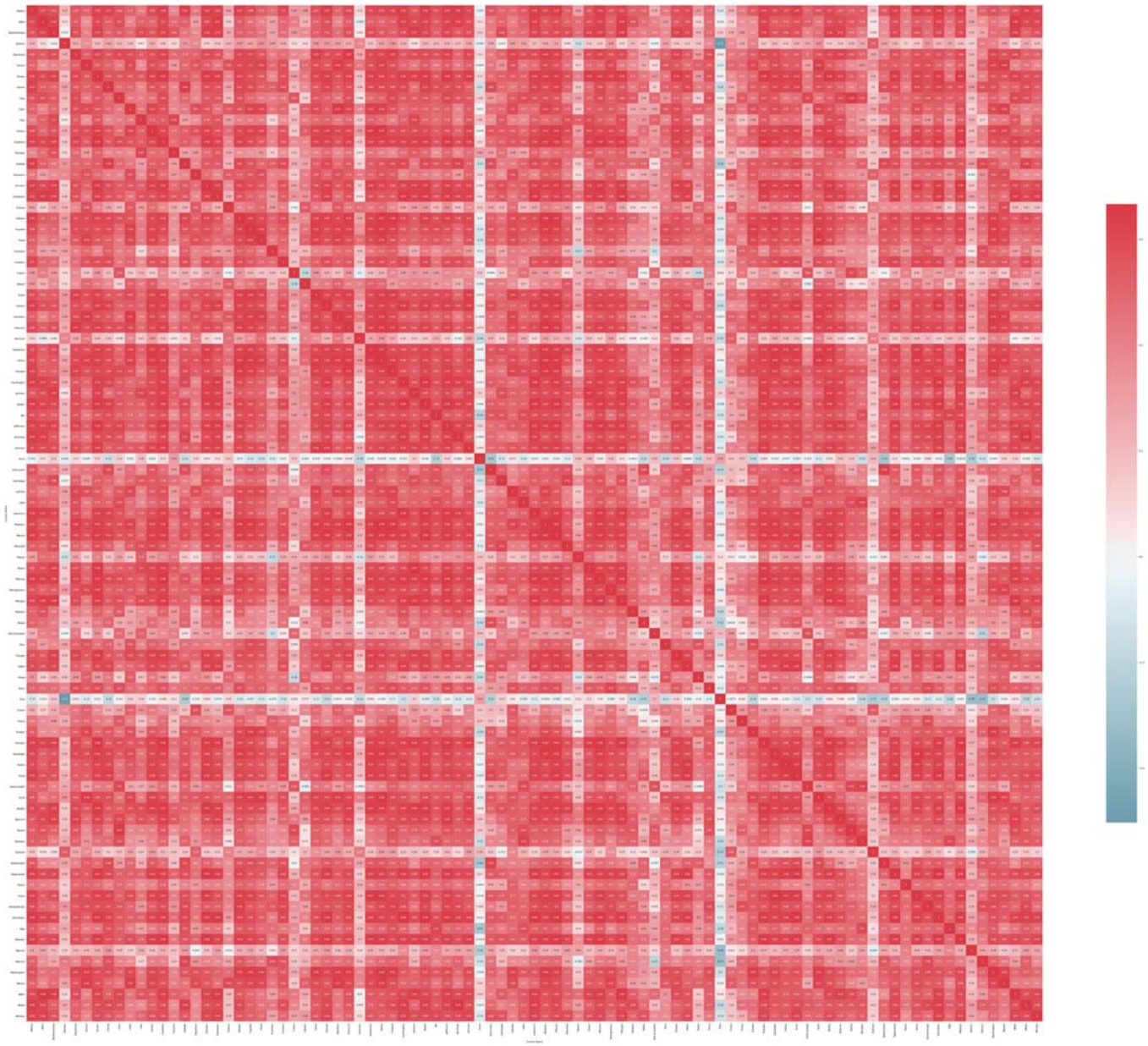
In [40]:
```
draw_heatmap_corr(create_pivot_df(df_grouped,'County Name','Year', 'Other Drugs'), 93,93
)
```



In [41]:
```
draw_heatmap_corr(create_pivot_df(df_grouped,'County Name','Year', 'Heroin'), 93,93)
```



In [42]:
```
draw_heatmap_corr(create_pivot_df(df_grouped,'County Name','Year', 'Methamphetamine'), 93,93)
```



Suplimentary analisys: PCA

The principal component analysis (PCA) (http://setosa.io/ev/principal-component-analysis/) is the most basic dimensionality reduction method. To run the PCA we want to isolate only the numerical columns.

In [43]:
```
from sklearn.decomposition import PCA
pca = PCA()
```

Now I ran `fit()` method to identify principal components.

In [44]:
```
pca_df_fitted = pca.fit(df_only_features)
print(pca.components_)
```

```
[[-3.85284978e-01  1.74017452e-01 -1.44132829e-01  5.72548369e-01
  -2.69179141e-02 -9.45944737e-04  4.88667622e-01  2.68324235e-02
   4.82126552e-01]
 [-4.18457857e-01 -5.05321504e-03  9.00800758e-01 -1.84296925e-02
  -5.53315489e-03  1.53651275e-03 -2.34787354e-02  1.09207683e-01
  -2.39850374e-02]
 [ 7.11756112e-01 -2.43212736e-01  3.74058306e-01  9.21887106e-02
  -1.39977393e-01 -4.67858705e-02  3.29558300e-01 -2.17148697e-01
   3.29070826e-01]
 [ 6.54654403e-02  2.33536949e-01  7.81258265e-02  1.44714164e-01
   6.66987977e-01 -5.86190546e-01 -6.12216977e-02 -3.43352728e-01
  -6.32252001e-02]
 [ 7.12073209e-02  4.07014556e-01  9.88869180e-02  2.09251566e-01
   5.82755415e-02  6.54588384e-01 -1.35008442e-01 -5.53985972e-01
  -1.36726374e-01]
 [-2.01654654e-01 -5.55680657e-01 -7.19214685e-02 -2.79439087e-01
   5.74989672e-01  3.75006655e-01  1.92924758e-01 -1.44138780e-01
   1.95080605e-01]
 [ 3.46163005e-01  2.35182719e-01  8.23741063e-02  1.93313001e-01
   4.47977083e-01  2.91666275e-01 -1.35950867e-02  7.03277331e-01
  -1.29745391e-02]
 [ 3.90868442e-04  5.68343537e-01  4.84514642e-03 -6.90996902e-01
   9.17264956e-03 -1.72477268e-03  2.33910579e-01  5.52881945e-03
   3.80333325e-01]
 [ 1.39976189e-03  5.68623039e-02  4.14804014e-04 -7.54235243e-02
   1.18822447e-03  5.79425375e-04  7.33450549e-01  1.74171781e-03
  -6.73142890e-01]]
```

'-Other Drugs', Marijuana,Opioid+Alcohol, Marijuana+Cocaine, Marijuana+Alcohol, -Marijuana-Heroin, Marijuana+Methamphetamine

In [45]:
```
print(pca.explained_variance_)
```

```
[0.52332706 0.25520259 0.20435398 0.13614426 0.12185475 0.1103172
 0.10171866 0.00989851 0.00189358]
```

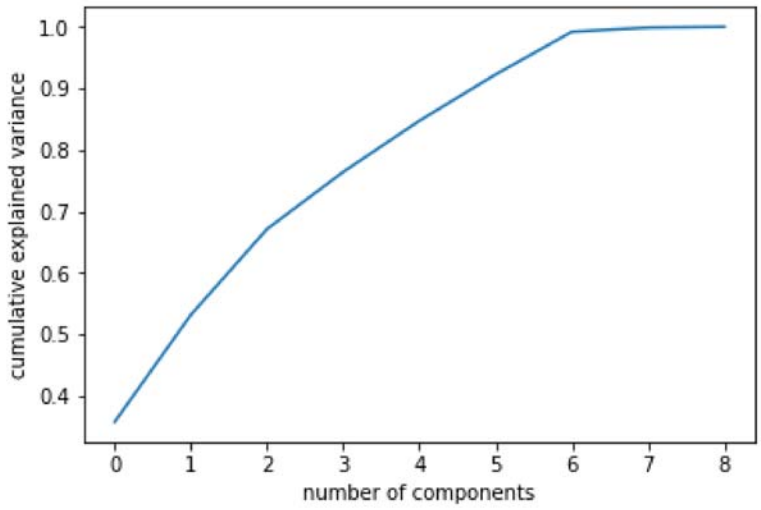
In [46]:
```
pca_df_fitted.explained_variance_ratio_
```

Out[46]:
```
array([0.35729042, 0.17423414, 0.13951833, 0.0929496 , 0.08319374,
       0.07531672, 0.06944625, 0.006758  , 0.0012928 ])
```

In [47]:
```
pca.components_.shape
```

Out[47]:
```
(9, 9)
```

In [48]:
```
plt.plot(np.cumsum(pca.explained_variance_ratio_))
plt.xlabel('number of components')
plt.ylabel('cumulative explained variance');
```



The first six components capture more than 95% of the variance in original dataset. This means that the PCA is not very effective on this dataset and six components will provide approximation for the rest of the dimensions.