Practical 9: Grouping Data

Aggregation, Classification & Clustering

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A common challenge in data analysis is how to group observations in a data set together in a way that allows for generalisation: *this* group of observations are similar to one another, *that* group is dissimilar to this group. Sometimes we have a *label* that we can use as part of the process (in which case we're doing **classification**), and somtimes we don't (in which case we're doing **clustering**). But what defines similarity and difference? There is no *one* answer to that question and so there are many different ways to cluster or classify data, each of which has strengths and weaknesses that make them more, or less, appropriate in different contexts.



🔗 Connections: This practical pulls together many topics covered in other modules, and many of t

1 Preamble

```
import warnings # This suppresses some meaningless errors from Seaborn and Pandas
warnings.simplefilter(action='ignore', category=FutureWarning)
```

```
import numpy as np
import pandas as pd
import geopandas as gpd
import seaborn as sns
import matplotlib.pyplot as plt
import matplotlib.cm as cm
import matplotlib as mpl
import re
import os
```

from matplotlib.colors import ListedColormap

```
# All of these are potentially useful, though
# not all have been used in this practical --
# I'd suggest exploring the use of different
# Scalers/Transformers as well as clustering
# algorithms...
```

from sklearn.neighbors import NearestNeighbors

from sklearn.decomposition import PCA

from sklearn.preprocessing import MinMaxScaler, StandardScaler, RobustScaler, PowerTransformer

from sklearn.cluster import KMeans, DBSCAN, OPTICS

 $from \ sklearn. ensemble \ import \ Random Forest Classifier$

 $from \ sklearn.metrics \ import \ silhouette_samples, \ silhouette_score$

```
import random
```

```
random.seed(42) # For reproducibility

np.random.seed(42) # For reproducibility

# Make numeric display a bit neater

pd.set_option('display.float_format', lambda x: '{:,.2f}'.format(x))
```

1.0.1 Initialise the Scaler(s)

Remember that you can set up the sklearn transformers in advance, and then fit them before transform-ing them.

```
mms = MinMaxScaler(feature_range=(-1,1))
stds = StandardScaler()
rbs = RobustScaler()
pts = PowerTransformer()
```

1.0.2 Set Up Plotting Functions

Note

🔗 Connections: Here's an example of how you can use a function to do something a little more co

```
def plt_ldn(w, b):
  Creates a new figure of a standard size with the
  water (w) and boundary (b) layers set up for easy
  plotting. Right now this function assumes that you're
 looking at London, but you could parameterise it in
  other ways ot allow it to work for other areas.
 w: a water layer for London
  b: a borough (or other) boundary layer for London
 fig, ax = plt.subplots(1, figsize=(14, 12))
 w.plot(ax=ax, color='#79aef5', zorder=2)
  b.plot(ax=ax, edgecolor='#cc2d2d', facecolor='None', zorder=3)
  ax.set_xlim([502000,563000])
  ax.set ylim([155000,201500])
  ax.spines['top'].set_visible(False)
  ax.spines['right'].set_visible(False)
  ax.spines['bottom'].set_visible(False)
  ax.spines['left'].set visible(False)
  return fig, ax
################################
# These may no longer be relevant because of changes to geopandas API
def default_cmap(n, outliers=False):
 cmap = mpl.cm.get_cmap('viridis_r', n)
  colors = cmap(np.linspace(0,1,n))
 if outliers:
    gray = np.array([225/256, 225/256, 225/256, 1])
    colors = np.insert(colors, 0, gray, axis=0)
 return ListedColormap(colors)
# mappable = ax.collections[-1] if you add the geopandas
# plot last.
def add_colorbar(mappable, ax, cmap, norm, breaks, outliers=False):
  cb = fig.colorbar(mappable, ax=ax, cmap=cmap, norm=norm,
           boundaries=breaks,
          extend=('min' if outliers else 'neither'),
          spacing='uniform',
          orientation='horizontal',
          fraction=0.05, shrink=0.5, pad=0.05)
  cb.set_label("Cluster Number")
```

1.0.3 Set up Caching Function

```
import os
from requests import get
from urllib.parse import urlparse
```

```
def cache_data(src:str, dest:str) -> str:
    """Downloads and caches a remote file locally.
```

The function sits between the 'read' step of a pandas or geopandas data frame and downloading the file from a remote location. The idea is that it will save it locally so that you don't need to remember to do so yourself. Subsequent re-reads of the file will return instantly rather than downloading the entire file for a second or n-th itme.

```
Parameters
_____
src:str
  The remote *source* for the file, any valid URL should work.
  The *destination* location to save the downloaded file.
Returns
  A string representing the local location of the file.
url = urlparse(src) # We assume that this is some kind of valid URL
fn = os.path.split(url.path)[-1] # Extract the filename
dfn = os.path.join(dest,fn) # Destination filename
# Check if dest+filename does *not* exist --
# that would mean we have to download it!
if not os.path.isfile(dfn) or os.path.getsize(dfn) < 1:</pre>
  print(f"{dfn} not found, downloading!")
  # Convert the path back into a list (without)
  # the filename -- we need to check that directories
  # exist first.
  path = os.path.split(dest)
  # Create any missing directories in dest(ination) path
  # -- os.path.join is the reverse of split (as you saw above)
  # but it doesn't work with lists... so I had to google how
  # to use the 'splat' operator! os.makedirs creates missing
  # directories in a path automatically.
  if len(path) >= 1 and path[0] != ":
    os.makedirs(os.path.join(*path), exist ok=True)
```

```
# Download and write the file
with open(dfn, "wb") as file:
response = get(src)
file.write(response.content)

print('Done downloading...')

else:
print(f"Found {dfn} locally!")

return dfn
```

2 Load Data

2.1 London Data Layers

```
   Difficulty: Low.
```

```
spath = 'https://github.com/jreades/fsds/blob/master/data/src/' # source path
ddir = os.path.join('data', 'geo') # destination directory
water = gpd.read_file( cache_data(spath+'Water.gpkg?raw=true', ddir) )
boros = gpd.read_file( cache_data(spath+'Boroughs.gpkg?raw=true', ddir) )
green = gpd.read_file( cache_data(spath+'Greenspace.gpkg?raw=true', ddir) )
msoas = gpd.read_file( cache_data('http://orca.casa.ucl.ac.uk/~jreades/data/MSOA-2011.gpkg', ddir) )
msoas = msoas.to_crs(epsg=27700)

# I don't use this in this practical, but it's a
# really useful data set that gives you 'names'
# for MSOAs that broadly correspond to what most
# Londoners would think of as a 'neighbourhood'.
msoa_nms = gpd.read_file( cache_data('http://orca.casa.ucl.ac.uk/~jreades/data/MSOA-2011-Names.gpkg', ddir) )
msoa_nms = msoa_nms.to_crs(epsg=27700)
print("Done.")
```

2.2 Reduced Dimensionality MSOA Data



You should have this locally from last week, but just in case...

```
host = 'http://orca.casa.ucl.ac.uk'
path = '~jreades/data'
rddf = gpd.read_parquet( cache_data(f'{host}/{path}/Reduced_Dimension_Data.geoparquet', ddir) )
print(f"Data frame is {rddf.shape[0]:,} x {rddf.shape[1]}")
```

You should have: Data frame is 983 x 93.

And below you should see both the components and the dimensions from last week's processing.

rddf.iloc[0:3, -7:]

I get the results below, but note that the **Dimension values** may be slightly different:

| Co | Component Component | | | Dimension Dimension | | | |
|-----------|---------------------|------|-------|--------------------------------------|------|------|---------------------------------------|
| | 5 | 6 | 7 | Borough | n 1 | 2 | Subregion |
| E02000001 | 1.44 | 3.95 | -1.52 | City of Lon- don | 7.74 | 3.36 | Inner West |
| E02000002 | 2 -0.28 | 0.89 | 0.26 | Barking and Da- gen- ham | 2.04 | 7.59 | Outer East and North East |
| E02000003 | 3 -0.11 | 1.12 | 0.83 | Barking and Da- gen- ham | 2.20 | 6.87 | Outer East and North East |

2.3 Listings Data



? Difficulty: Low.

Let's also get the listings data from a few weeks back:

listings = gpd.read_parquet(cache_data(f{host}/{path}/2023-09-06-listings.geoparquet', ddir)) listings = listings.to_crs(epsg=27700) print(f"Data frame is {listings.shape[0]:,} x {listings.shape[1]}")

You should have: Data frame is 85,134 x 31.

And a quick plot of the price to check:

listings.plot(???, cmap='plasma', scheme='quantiles', k=10, markersize=.5, alpha=0.15, figsize=(10,7));

3 Aggregate Listings by MSOA

3.1 Join Listings to MSOA



🛕 Difficulty: Medium-to-hard.

First, let's link all this using the MSOA Geography that we created last week and a mix or merge and sjoin!

```
Note
  **🔗 Connections**: Notice a few things going on here! We are calling `gpd.sjoin` because pandas (`p
# Before the spatial join
listings.columns
msoa_listings = gpd.sjoin(???, msoas.drop(
            columns=['MSOA11NM', 'LAD11CD', 'LAD11NM', 'RGN11CD', 'RGN11NM',
                 'USUALRES', 'HHOLDRES', 'COMESTRES', 'POPDEN', 'HHOLDS',
                 'AVHHOLDSZ']), predicate='???').drop(
            columns=['latitude','longitude','index_right']
        )
# All we've added is the MSOA11CD
msoa_listings.columns
All being well you should now have:
Index(['listing_url', 'last_scraped', 'name', 'description', 'host_id',
   'host_name', 'host_since', 'host_location', 'host_is_superhost',
   'host_listings_count', 'host_total_listings_count',
   'host_verifications', 'property_type', 'room_type', 'accommodates',
   'bathrooms_text', 'bedrooms', 'beds', 'amenities', 'price',
   'minimum_nights', 'maximum_nights', 'availability_365',
```

3.2 Price by MSOA

dtype='object')



🛕 Difficulty: Medium.

'number_of_reviews', 'first_review', 'last_review',

'review_scores_rating', 'reviews_per_month', 'geometry', 'MSOA11CD'],

Let's calculate the median price by MSOA... Notice that we have to specify the column we want after the groupby so the we don't get the median of every column returned

Note

🔗 Connections: I find `groupby` to be a complex operation and often need a couple of gos before

```
# *m*soa *|*istings *g*rouped by *p*rice
mlgp = msoa_listings.groupby('???')['price'].agg('???')
mlgp.head()
```

You should get something like:

```
MSOA11CD

E02000001 170.00

E02000002 97.00

E02000003 80.00

E02000004 54.00

E02000005 100.00

Name: price, dtype: float64
```

3.3 Room Type by MSOA

```
▲ Difficulty: Medium.
```

Now let's calculate the count of room types by MSOA and compare the effects of reset_index on the outputs below. And notice too that we can assign the aggregated value to a column name!

```
# *m*soa *l*istings *g*rouped *c*ount
mlgc = msoa_listings.groupby(['???','???'], observed=False).listing_url.agg(Count='???')
mlgc.head()
```

You should get something resembling this:

| MSOA11CD | room_type | Count |
|-----------|-----------------|-------|
| E0200001 | Entire home/apt | 466 |
| | Hotel room | 0 |
| | Private room | 61 |
| | Shared room | 1 |
| E02000002 | Entire home/apt | 4 |

*m*soa *l*istings *g*rouped *c*ount *r*eset index
mlgcr = msoa_listings.groupby(['???','???'], observed=False).listing_url.agg(Count='???').reset_index() # msoa listing
mlgcr.head()

You should get something like:

| MSOA11CD | room_type | Count |
|-----------|--|---|
| E02000001 | Entire home/apt | 466 |
| E02000001 | Hotel room | 0 |
| E02000001 | Private room | 61 |
| E02000001 | Shared room | 1 |
| E02000002 | Entire home/apt | 4 |
| | E02000001 E02000001 E02000001 E02000001 | E02000001 Entire home/apt E02000001 Hotel room E02000001 Private room E02000001 Shared room |

3.4 Price by Room Type



Difficulty: Hard.

But perhaps median price/room type would make more sense? And do we want to retain values where there are no listings? For example, there are no hotel rooms listed for E02000001, how do we ensure that these *NAs are dropped?*

```
# *m*soa *l*istings *g*rouped *r*oom *p*rice
mlgrp = msoa_listings.???(???, observed=True
           )['price'].agg('???').reset_index()
mlgrp.head()
```

You should get something like:

| | MSOA11CD | room type | price |
|---|-----------|-----------------|--------|
| 0 | E02000001 | Entire home/apt | 177.00 |
| 2 | E02000001 | Private room | 100.00 |
| 3 | E02000001 | Shared room | 120.00 |
| 4 | E02000002 | Entire home/apt | 117.00 |
| 6 | E02000002 | Private room | 42.00 |

3.5 Explore Outlier Per-MSOA Prices



🛕 Difficulty: Medium.

Are there MSOAs what look like they might contain erroneous data?

3.5.1 Plot MSOA Median Prices

mlgp.hist(bins=200);

3.5.2 Examine Listings from High-Priced MSOAs

Careful, this is showing the listings from MSOAs whose median price is above \$300/night:

```
msoa_listings[
  msoa_listings.MSOA11CD.isin(mlgp[mlgp > 300].index)
].sort_values(by='price', ascending=False).head(7)[
  ['price','room type','name','description']
1
```

Some of these look legi (4, 5, and... 8 bedroom 'villas'?), though not every one...

And how about these?

```
msoa_listings[
   (msoa_listings.MSOA11CD.isin(mlgp[mlgp > 300].index)) & (msoa_listings.room_type!='Entire home/apt')
].sort_values(by='price', ascending=False).head(7)[
   ['price','room_type','property_type','name','description']
]
```

If we wanted to be rigorous then we'd have to investigate further: properties in Mayfair and Westminster *are* going to be expensive, but are these plausible nightly prices? In some cases, yes. In others...

```
msoa_listings[
   (msoa_listings.MSOA11CD.isin(mlgp[mlgp < 100].index)) & (msoa_listings.room_type!='Entire home/apt')
].sort_values(by='price', ascending=False).head(7)[
   ['price','room_type','name','description']
]
```

On the whole, let's take a *guess* that there are a small number of implausibly high prices for individual units that aren't in very expensive neighbourhoods and that these are either erroneous/deliberately incorrect, or represent a price that is not per-night.

Note

🔗 Connections: What's the right answer here? There isn't one. You could probably spend _months

3.5.3 Filter Unlikely Listings

Difficulty: Hard.

See if you can filter out these less likely listings on the following criteria:

- 1. Listings are priced above \$300/night AND
- 2. Room type is not 'Entire home/apt' AND
- 3. Listings do *not* contain the words: suite, luxury, loft, stunning, prime, historic, or deluxe.

I found 901 rows to drop this way.

3.5.4 Plot Unlikely Listings

Here we use the plt_ldn function – notice how it's designed to return f,ax in the same way that plt.subplots (which we're already familiar with) does!

```
f,ax = plt_ldn(???, ???)
to_drop.plot(column='price', markersize=10, alpha=0.7, cmap='viridis', ax=ax);
```

3.5.5 ... And Drop

Some might be legitimate, but I'm feeling broadly ok with the remainder.

```
cleaned = msoa_listings.drop(index=to_drop.???)
print(f"Cleaned data has {cleaned.shape[0]:,} rows.")
```

After this I had 84,308 rows.

I would normally, at this point, spend quite a bit of time validating this cleaning approach, but right now we're going to take a rough-and-ready approach.

3.5.6 Questions

- What data type did Task 2.2 return?
- What is the function of reset_index() in Task 2.3 and when might you choose to reset (or not)?

4 Pivot Tables & 'Wide Data'

The group_by operation is *one* way to organise and aggregate our data, but pivot tables are a *second* common way to achieve this. We typically use a pivot table to go from long to wide data frames – it's often seen as one of Excel's main benefits, but Pandas can do that too!

Note

🔗 Connections: Notice that a pivot table is just a different kind of aggregation. Principally, it's about

4.1 Create Pivot Table



Difficulty: Hard.

We can make use of the pivot table function to generate counts by MSOA in a 'wide' format.

```
pivot = cleaned.groupby(
        ['MSOA11CD','room_type'], observed=False
    ).listing_url.agg(Count='count').reset_index().pivot(
         index='???', columns=['???'], values=['???'])
pivot.head(3)
```

The formatting will look a tiny bit different, but you should get something like this:

| | | | | | Count |
|----------------|------------|------------|---------|--------|-------|
| room_typEntire | e home/apt | Hotel room | Private | Shared | |
| | | | room | room | |
| MSOA11CD | | | | | |
| E0200001 | 466 | 0 | 55 | 1 | |
| E02000002 | 4 | 0 | 2 | 0 | |
| E0200003 | 12 | 0 | 13 | 0 | |

4.2 Check Counts



Difficulty: Low.

pivot.sum()

Just to reassure you that the pivot results 'make sense':

```
print(cleaned[cleaned.room_type=='Entire home/apt'].listing_url.count())
print(cleaned[cleaned.room_type=='Private room'].listing_url.count())
```

4.3 Tidy & Normalise



Difficulty: Low.

My instinct at this point is that, looking at the pivot table, we see quite different levels of Airbnb penetration and it is hard to know how handle this difference: share would be unstable because of the low counts in some places and high counts in others; a derived variable that tells us something about density or mix could be interesting (e.g. HHI or LQ) but wouldn't quite capture the pattern of mixing.

4.3.1 Tidy

Personally, based on the room type counts above I think we can drop Hotel Rooms and Shared Rooms from this since the other two categories are so dominant.

```
# Flatten the column index
pivot.columns = ['Entire home/apt','Hotel room','Private room','Shared room']
# Drop the columns
pivot.drop(???, inplace=True)
pivot.head()
```

You should have only the Entire home/apt and Private room columns now.

4.3.2 Normalise

```
pivot_norm = pd.DataFrame(index=pivot.index)
for c in pivot.columns.to_list():
  # Power Transform
  pivot norm[c] = pts.???(pivot[c].to numpy().reshape(???,???))
```

pivot norm.head()

You should have something like:

| | Entire home/apt | Private room | |
|-----------|-----------------|--------------|--|
| MSOA11CD | | | |
| E0200001 | 2.20 | 1.06 | |
| E02000002 | -1.29 | -1.85 | |

4.3.3 Plot

pnm = pd.merge(msoas.set index('MSOA11CD'), pivot norm, left index=True, right index=True) pnm.plot(column='Entire home/apt', cmap='viridis', edgecolor='none', legend=True, figsize=(12,8));

4.4 PCA



🛕 Difficulty: Moderate, though you might find the questions hard.

You can merge the output of this next step back on to the rddf data frame as part of a clustering process, though we'd really want to do some more thinking about what this data means and what transformations we'd need to do in order to make them meaningful.

For instance, if we went back to last week's code, we could have appended this InsideAirbnb data before doing the dimensionality reduction, or we could apply it now to create a new measure that could be used as a separate part of the clustering process together with the reduced dimensionality of the demographic data.

4.4.1 Perform Reduction

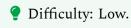
```
pcomp = PCA(n components=???, random state=42)
rd = pcomp.???(pivot_norm)
print(f"The explained variance of each component is: {', '.join([f'{x*100:.2f}%' for x in pcomp.explained_variance_ra
Take the first component and convert to a series to enable the merge:
airbnb pca = pd.DataFrame(
        {'Airbnb Component 1': mms.fit transform(rd[:,1].reshape(-1,1)).reshape(1,-1)[0]},
        index=pivot.index)
airbnb_pca.head()
You should have something like: | | Airbnb Component 1 | | :-- | ---: | | MSOA11CD |
| | E02000001 | 0.47 | | E02000002 | 0.19
pcanm = pd.merge(msoas.set index('MSOA11CD'), airbnb pca, left index=True, right index=True)
pcanm.plot(column='Airbnb Component 1', cmap='viridis', edgecolor='none', legend=True, figsize=(12,8));
4.4.2 Write to Data Frame
# Result Set from merge
rs = pd.merge(rddf, airbnb_pca, left_index=True, right_index=True)
Grab the PCA, UMAP, and Airbnb outputs for clustering and append rescaled price:
# Merge the reducded dimensionality data frame with the PCA-reduced Airbnb data
# to create the *cl*uster *d*ata *f*rame
cldf = pd.merge(rddf.loc[:,'Component 1':], airbnb_pca,
        left_index=True, right_index=True)
# Append median price from cleaned listings grouped by MSOA too!
s1 = cleaned.groupby(by='MSOA11CD').price.agg('median')
cldf['median_price'] = pd.Series(np.squeeze(mms.fit_transform(s1.values.reshape(-1,1))), index=s1.index)
# Append mean price from cleaned listings grouped by MSOA too!
s2 = cleaned.groupby(by='MSOA11CD').price.agg('mean')
cldf['mean_price'] = pd.Series(np.squeeze(mms.fit_transform(s2.values.reshape(-1,1))), index=s2.index)
cldf.drop(columns=['Subregion','Borough'], inplace=True)
cldf.head()
```

4.4.3 Questions

 Have a think about why you might want to keep the Airbnb data separate from the MSOA data when doing PCA (or any other kind of dimensionality reduction)! • Why *might* it be interesting to add *both* mean and median MSOA prices to the clustering process? Here's a hint (but it's *very* subtle): sns.jointplot(x=s1, y=s2, s=15, alpha=0.6)

5 First K-Means Clustering

5.1 Perform Clustering



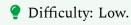
```
c_nm = 'KMeans' # Clustering name
k pref = ??? # Number of clusters
```

kmeans = KMeans(n_clusters=k_pref, n_init=25, random_state=42).fit(cldf.drop(columns=['Dimension 1','Dimension 1',

Here are the results:

print(kmeans.labels_) # The results

5.2 Save Clusters to Data Frame



5.2.1 Write Series and Assign

Now capture the labels (i.e. clusters) and write them to a data series that we store on the result set df (rs):

rs[c_nm] = pd.Series(kmeans.labels_, index=cldf.index)

5.2.2 Histogram of Cluster Members

How are the clusters distributed?

sns.histplot(data=???, x=c_nm, bins=k_pref);

5.2.3 Map Clusters

And here's a map!

```
fig, ax = plt_ldn(water, boros)
fig.suptitle(f"{c_nm} Results (k={k_pref})", fontsize=20, y=0.92)
rs.plot(column=???, ax=ax, linewidth=0, zorder=0, categorical=???, legend=True);
```

5.2.4 Questions

- What critical assumption did we make when running this analysis?
- Why did I not use the UMAP dimensions here?
- Why do we have the c nm='kMeans' when we know what kind of clustering we're doing?
- Does this look like a good clustering?

6 Second K-Means Clustering

6.1 What's the 'Right' Number of Clusters?



Difficulty: Moderate.

There's more than one way to find the 'right' number of clusters. In Singleton's Geocomputation chapter they use WCSS to pick the 'optimal' number of clusters. The idea is that you plot the average WCSS for each number of possible clusters in the range of interest (2...n) and then look for a 'knee' (i.e. kink) in the curve. The principle of this approach is that you look for the point where there is declining benefit from adding more clusters. The problem is that there is always some benefit to adding more clusters (the perfect clustering is k==n), so you don't always see a knee.

Another way to try to make the process of selecting the number of clusters a little less arbitrary is called the silhouette plot and (like WCSS) it allows us to evaluate the 'quality' of the clustering outcome by examining the distance between each observation and the rest of the cluster. In this case it's based on Partitioning Around the Medoid (PAM).

Either way, to evaluate this in a systematic way, we want to do multiple k-means clusterings for multiple values of k and then we can look at which gives the best results...

kcldf = cldf.drop(columns=['Dimension 1','Dimension 2'])

6.1.1 Repeated Clustering

Let's try clustering across a wider range. Because we repeatedly re-run the clustering code (unlike with Hierarchical Clustering) this can take a few minutes. I got nearly 5 minutes on a M2 Mac.

```
%%time
# Adapted from: http://scikit-learn.org/stable/auto examples/cluster/plot kmeans silhouette analysis.html
x = []
y = []
# For resolutions of 'k' in the range 2..40
for k in range(2,41):
  ###############
  # Do the clustering using the main columns
  kmeans = KMeans(n_clusters=k, n_init=25, random_state=42).fit(kcldf)
  # Calculate the overall silhouette score
  silhouette_avg = silhouette_score(kcldf, kmeans.labels_)
  y.append(k)
  x.append(silhouette avg)
  print('.', end=")
6.1.2 Plot Silhouette Scores
print(f"Largest silhouette score was {max(x):6.4f} for k={y[x.index(max(x))]}")
plt.plot(y, x)
plt.gca().xaxis.grid(True);
plt.gcf().suptitle("Average Silhouette Scores");
```

Warning

⚠ Note: Had we used the UMAP dimensions here you'd likely see more instability in the silhouette plants.

We can use the largest average silhouette score to determine the 'natural' number of clusters in the data, but that that's only if we don't have any kind of underlying theory, other empirical evidence, or even just a reason for choosing a different value... Again, we're now getting in areas where your judgement and your ability to communicate your rationale to readers is the key thing.

6.2 Final Clustering



Difficulty: Low.

So although we should probably pick the largest silhouette scores, that's k=3 which kind of defeats the purpose of clustering in the first place. In the absence of a compelling reason to pick 2 or 3 clusters, let's have a closer look at the next maximum silhouetted score:

6.2.1 Perform Clustering k pref=??? ############## # Do the clustering using the main columns kmeans = KMeans(n clusters=k pref, n init=25, random state=42).fit(kcldf) # Convert to a series s = pd.Series(kmeans.labels_, index=kcldf.index, name=c_nm) # We do this for plotting $rs[c_nm] = s$ # Calculate the overall silhouette score silhouette_avg = silhouette_score(kcldf, kmeans.labels_) # Calculate the silhouette values sample_silhouette_values = silhouette_samples(kcldf, kmeans.labels_) 6.2.2 Plot Diagnostics ############## # Create a subplot with 1 row and 2 columns fig, (ax1, ax2) = plt.subplots(1, 2)fig.set_size_inches(9, 5) # The 1st subplot is the silhouette plot # The silhouette coefficient can range from -1, 1 ax1.set_xlim([-1.0, 1.0]) # Changed from -0.1, 1 # The (n_clusters+1)*10 is for inserting blank space between silhouette # plots of individual clusters, to demarcate them clearly. $ax1.set_ylim([0, kcldf.shape[0] + (k_pref + 1) * 10])$ $y_lower = 10$ # For each of the clusters... for i in range(k_pref):

```
# Aggregate the silhouette scores for samples belonging to
# cluster i, and sort them
ith cluster silhouette values = \
  sample_silhouette_values[kmeans.labels_ == i]
ith_cluster_silhouette_values.sort()
size cluster i = ith cluster silhouette values.shape[0]
y_upper = y_lower + size_cluster_i
# Set the color ramp
color = plt.cm.Spectral(i/k pref)
ax1.fill_betweenx(np.arange(y_lower, y_upper),
           0, ith_cluster_silhouette_values,
           facecolor=color, edgecolor=color, alpha=0.7)
# Label the silhouette plots with their cluster numbers at the middle
ax1.text(-0.05, y_lower + 0.5 * size_cluster_i, str(i))
# Compute the new y lower for next plot
y_lower = y_upper + 10 # 10 for the 0 samples
ax1.set_title("The silhouette plot for the clusters.")
ax1.set xlabel("The silhouette coefficient values")
ax1.set_ylabel("Cluster label")
# The vertical line for average silhouette score of all the values
ax1.axvline(x=silhouette_avg, color="red", linestyle="--", linewidth=0.5)
ax1.set yticks([]) # Clear the yaxis labels / ticks
ax1.set_xticks(np.arange(-1.0, 1.1, 0.2)) # Was: [-0.1, 0, 0.2, 0.4, 0.6, 0.8, 1]
# 2nd Plot showing the actual clusters formed --
# we can only do this for the first two dimensions
# so we may not see fully what is causing the
# resulting assignment
colors = plt.cm.Spectral(kmeans.labels .astype(float) / k pref)
ax2.scatter(kcldf[kcldf.columns[0]], kcldf[kcldf.columns[1]],
       marker='.', s=30, lw=0, alpha=0.7, c=colors)
# Labeling the clusters
centers = kmeans.cluster centers
# Draw white circles at cluster centers
ax2.scatter(centers[:, 0], centers[:, 1],
      marker='o', c="white", alpha=1, s=200)
for i, c in enumerate(centers):
  ax2.scatter(c[0], c[1], marker='$%d$' % i, alpha=1, s=50)
ax2.set_title("Visualization of the clustered data")
ax2.set_xlabel("Feature space for the 1st feature")
```

```
ax2.set_ylabel("Feature space for the 2nd feature")
plt.suptitle(("Silhouette results for KMeans clustering "
        "with %d clusters" % k pref),
        fontsize=14, fontweight='bold')
```

plt.show()



Warning

⚠ Stop: Make sure that you understand how the silhouette plot and value work, and why your results

6.2.3 Map Clusters

```
fig, ax = plt_ldn(water, boros)
fig.suptitle(f"{c_nm} Results (k={k_pref})", fontsize=20, y=0.92)
rs.plot(column=c_nm, ax=ax, linewidth=0, zorder=0, categorical=True, legend=True);
```

6.3 'Representative' Centroids



A Difficulty: Moderate since, conceptually, there's a lot going on.

To get a sense of how these clusters differ we can try to extract 'representative' centroids (mid-points of the multi-dimensional cloud that constitutes a cluster). In the case of k-means this will work quite will since the clusters are explicitly built around mean centroids. There's also a k-medoids clustering approach built around the median centroid.

These are columns that we want to suppress from our sample:

```
to_suppress=['OBJECTID', 'BNG_E', 'BNG_N', 'LONG', 'LAT',
      'Shape__Are', 'Shape__Len', 'geometry', 'Component 1',
      'Component 2', 'Component 3', 'Component 4', 'Component 5',
      'Component 6', 'Component 7', 'Dimension 1', 'Dimension 2',
      'Airbnb Component 1']
```

Take a sample of the full range of numeric columns:

cols = random.sample(rs.select_dtypes(exclude='object').drop(columns=to_suppress).columns.to_list(), 12) print(cols)

Calculate the mean of these columns for each cluster:

```
# Empty data frame with the columns we'll need
centroids = pd.DataFrame(columns=cols)
# For each cluster...
for k in sorted(rs[c_nm].unique()):
  print(f"Processing cluster {k}")
```

```
# Select rows where the cluster name matches the cluster number
 clust = rs[rs[c_nm] == k]
  # Append the means to the centroids data frame
 centroids.loc[k] = clust[cols].mean()
centroids
centroids_long = pd.DataFrame(columns=['Variable','Cluster','Std. Value'])
for i in range(0,len(centroids.index)):
  row = centroids.iloc[i,:]
 for r in row.index:
    d = pd.DataFrame({'Variable':r, 'Cluster':i, 'Std. Value':row[r]}, index=[1])
    centroids_long = pd.concat([centroids_long, d], ignore_index=True)
g = sns.FacetGrid(centroids_long, col="Variable", col_wrap=3, height=3, aspect=1.5, margin_titles=True, sharey=Tru
g = g.map(plt.bar, "Cluster", "Std. Value")
  Note
```

🔗 Connections: The above centroid outputs are a way to think about how each cluster is 'loaded' o

7 DBSCAN

For what it's worth, I've had enormous trouble with DBSCAN and this kind of data. I don't think it deals very well with much more than three dimensions, so the flexbility to not have to specify the number of clusters is balanced with a density-based approach that is severely hampered by high-dimensional distance-inflation.

```
# Drop the PCA dimensions
cldf2 = cldf.loc[:,'Dimension 1':].copy()
for c in [x for x in cldf.columns.to_list() if x.startswith('Dimension ')]:
  cldf2[c] = pd.Series(np.squeeze(mms.fit_transform(cldf2[c].to_numpy().reshape(-1,1))), index=cldf2.index)
cldf2.head()
```

7.1 Work out the Neighbour Distance

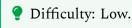


Difficulty: Moderate.

We normally look for some kind of 'knee' to set the distance.

```
nbrs = NearestNeighbors(n_neighbors=6).fit(cldf2)
distances, indices = nbrs.kneighbors(cldf2)
distances = np.sort(distances, axis=0)
distances = distances[:,1]
```

7.2 Derive Approximate Knee



from kneed import knee_locator

```
kn = knee locator.KneeLocator(np.arange(distances.shape[0]), distances, S=12,
                curve='convex', direction='increasing')
print(f"Knee detected at: {kn.knee}")
kn.plot_knee()
kn.plot_knee_normalized()
print(f"Best guess at epsilon for DBSCAN is {distances[kn.knee]:0.4f}")
```

7.3 Explore Epsilons



🛕 Difficulty: Moderate.

There are two values that need to be specified: eps and min_samples. Both seem to be set largely by trial and error, though we can use the above result as a target. It's easiest to set min_samples first since that sets a floor for your cluster size and then eps is basically a distance metric that governs how far away something can be from a cluster and still be considered part of that cluster.

7.3.1 Iterate Over Range



Caution

⚠ Warning: Depending on the data volume, this next step may take quite a lot of time since we are it

%%time

```
c_nm = 'DBSCAN'
# Make numeric display a bit neater
pd.set option('display.float format', lambda x: '{:,.4f}'.format(x))
el = []
max clusters = 10
cluster_count = 1
iters = 0
for e in np.arange(0.025, 0.76, 0.01): # <- You might want to adjust these!
  if iters % 25==0: print(f"{iters} epsilons explored.")
```

```
# Run the clustering
  dbs = DBSCAN(eps=e, min_samples=cldf2.shape[1]+1).fit(cldf2)
  # See how we did
  s = pd.Series(dbs.labels_, index=cldf2.index, name=c_nm)
  row = [e]
  data = s.value_counts()
  for c in range(-1, max_clusters+1):
    try:
      if np.isnan(data[c]):
        row.append(None)
      else:
        row.append(data[c])
    except KeyError:
      row.append(None)
  el.append(row)
  iters+=1
edf = pd.DataFrame(el, columns=['Epsilon']+["Cluster " + str(x) for x in list(range(-1,max_clusters+1))])
# Make numeric display a bit neater
pd.set_option('display.float_format', lambda x: '{:,.2f}'.format(x))
print("Done.")
7.3.2 Examine Clusters
edf.head() # Notice the -1 cluster for small epsilons
epsilon_long = pd.DataFrame(columns=['Epsilon','Cluster','Count'])
for i in range(0,len(edf.index)):
 row = edf.iloc[i,:]
 for c in range(1,len(edf.columns.values)):
    if row[c] != None and not np.isnan(row[c]):
      d = pd.DataFrame({'Epsilon':row[0], 'Cluster':f"Cluster {c-2}", 'Count':row[c]}, index=[1])
      epsilon_long = pd.concat([epsilon_long, d], ignore_index=True)
epsilon_long['Count'] = epsilon_long.Count.astype(float)
```

7.3.3 Plot Cluster Sizes

One of the really big problems with DBSCAN and this kind of data is that you have no *practical* way of specifying epsilon (whereas if you were doing walkability analysis then you could cluster on walking distance!). So you can look at the data (as above) to get a reasoanble value, but look what the output below shows about the stability of the clusters for different values of epsilon!

```
fig, ax = plt.subplots(figsize=(12,8))
sns.lineplot(data=epsilon_long, x='Epsilon', y='Count', hue='Cluster');
plt.vlines(x=distances[kn.knee], ymin=0, ymax=epsilon long.Count.max(), color=(1, .7, .7, .8), linestyles='dashed')
plt.gcf().suptitle(f"Cluster sizes for various realisations of Epsilon");
plt.tight_layout()
```

7.4 Final Clustering



Difficulty: Moderate.

###: Perform Clustering

Use the value from kneed...

```
dbs = DBSCAN(eps=distances[kn.knee], min_samples=cldf2.shape[1]+1).fit(cldf2.values)
s = pd.Series(dbs.labels , index=cldf2.index, name=c nm)
rs[c nm] = s
print(s.value_counts())
###: Map Clusters
fig, ax = plt_ldn(water, boros)
fig.suptitle(f"{c nm} Results", fontsize=20, y=0.92)
rs.plot(column=c nm, ax=ax, linewidth=0, zorder=0, legend=True, categorical=True);
```

7.4.1 'Representative' Centroids

```
to_suppress=['OBJECTID', 'BNG_E', 'BNG_N', 'LONG', 'LAT',
      'Shape_Are', 'Shape_Len', 'geometry', 'Component 1',
      'Component 2', 'Component 3', 'Component 4', 'Component 5',
      'Component 6', 'Component 7', 'Dimension 1', 'Dimension 2',
      'Airbnb Component 1']
```

Take a sample of the full range of numeric columns:

cols = random.sample(rs.select_dtypes(exclude='object').drop(columns=to_suppress).columns.to_list(), 12) print(cols)

Calculate the mean of these columns for each cluster:

```
# Empty data frame with the columns we'll need
centroids = pd.DataFrame(columns=cols)
# For each cluster...
for k in sorted(rs[c_nm].unique()):
  print(f"Processing cluster {k}")
  # Select rows where the cluster name matches the cluster number
  clust = rs[rs[c_nm]==k]
```

```
# Append the means to the centroids data frame
centroids.loc[k] = clust[cols].mean()

# Drop the unclustered records (-1)
centroids.drop(labels=[-1], axis=0, inplace=True)
centroids

centroids

centroids_long = pd.DataFrame(columns=['Variable','Cluster','Std. Value'])

for i in range(0,len(centroids.index)):
    row = centroids.iloc[i,:]
    for r in row.index:
        d = pd.DataFrame({'Variable':r, 'Cluster':i, 'Std. Value':row[r]}, index=[1])
        centroids_long = pd.concat([centroids_long, d], ignore_index=True)

g = sns.FacetGrid(centroids_long, col="Variable", col_wrap=3, height=3, aspect=1.5, margin_titles=True, sharey=True
g = g.map(plt.bar, "Cluster", "Std. Value")
```

8 Self-Organising Maps

SOMs offer a third type of clustering algorithm. They are a relatively 'simple' type of neural network in which the 'map' (of the SOM) adjusts to the data. We're not going to do this in *this* practical, but the main thing is that, unlike the above approaches, SOMs build a 2D map of a higher-dimensional space and use this as a mechanism for subsequently clustering the raw data. In this sense there is a conceptual link between SOMs and PCA or t-SNE or UMAP. Tehy are used quite a lot for text-clustering using keywords (where you have high-dimensionality).

There are a lot of SOM implementations in Python but the one I used to use, called SOMPY, appears to have been abandonned.

9 Classification

And now for something completely different! This is section is **completely optional**, but I thought that you might find it helpful to have a look at how *supervised* learning (classification) differs from *unsupervised* learning (clustering). Here we're going to perform a fairly straightforward classification: predicting the room_type for randomly-selected listings. Of course we know the true answer, but this is for demonstration purposes!

9.1 Additional Setup

Difficulty: Hard, as I've left out quite a bit of code.

9.1.1 Import Libraries

```
from sklearn.ensemble import RandomForestClassifier from sklearn.model_selection import train_test_split from sklearn.preprocessing import LabelEncoder from sklearn.metrics import confusion_matrix from sklearn.inspection import permutation importance
```

9.1.2 Set Up Data

I'm taking a fairly brutal approach here: anything that is not inherently numeric is gone (bye, bye, text), and I'm not bothering to convert implicitly numeric values either: dates could be converted to 'months since last review', for instance, while amenities could be One-Hot Encoded after some pruning of rare amenities. This leaves us with a much smaller number of columns to feed *in* to the classifier.

9.1.3 Remove NAs

Not all classifiers have this issue, but some will struggle to make predictions (or not be able to do so at all) if there are NAs in the data set. The classifier we're using can't deal with NAs, so we have to strip these out, but before we do let's check the effect:

```
classifier_in.isna().sum()
```

We can safely drop these now, and you should end up with about 54,000 rows to work with.

```
classifier_in = classifier_in.dropna(axis=0, how='any')
print(f"Now have {classifier_in.shape[0]:,} rows of data to work with (down from {cleaned.shape[0]:,}).")
print()
print(f"Classifier training columns: {', '.join(classifier_in.columns.to_list())}.")
classifier_in.head()
```

9.1.4 Remap Non-Numeric Columns

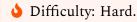
We do still have a couple of non-numeric columns to deal with: booleans and the thing we're actually trying to predict (the room type)!

```
classifier_in['host_is_superhost'] = classifier_in.host_is_superhost.replace({True:1, False:0}).astype('int')
le = LabelEncoder()
classifier_in['room_class'] = le.fit_transform(classifier_in.room_type)
```

A quick check: we should only have one type per class and vice versa.

classifier_in.groupby(by=['room_type','room_class']).host_id.agg('count').reset_index()

9.2 Random Forest Classification



We're going to use a Random Forest Classifier but the nice thing about sklearn is that you can quite easily swap in other classifiers if you'd like to explore further. This is one big advantage of Python over R in my book: whereas R tends to get new algorithms first, they are often implemented independently by many people and you can end up with incompatible data structures that require a lot of faff to reorganise for a different algorithm. Python is a bit more 'managed' and the dominance of numpy and sklearn and pandas means that people have an incentive to contribute to this library or, if it's genuinely novel, to create an implementation that works like it would if it were part of sklearn!

i Note

🔗 Connections: So here's an _actual_ Machine Learning implementation, but you'll have seen a lo

9.2.1 Train/Test Split

```
train, test = train_test_split(classifier_in, test_size=0.2, random_state=42)
print(f"Train contains {train.shape[0]:,} records.")
print(f"Test contains {test.shape[0]:,} records.")
y_train = train.room_class
X_train = train.drop(columns=['room_class','room_type'])
y test = test.room class
X_test = test.drop(columns=['room_class','room_type'])
9.2.2 Classifier Setup
rfc = RandomForestClassifier(
  max_depth=8,
  min_samples_split=7,
  n_jobs=4,
  random_state=42
)
9.2.3 Fit and Predict
rfc.fit(X_train, y_train)
y_hat = rfc.predict(X_test)
```

9.3 Validate



Difficulty: Hard.

9.3.1 Confusion Matrix

```
c_matrix = pd.DataFrame(confusion_matrix(y_test, y_hat))
c matrix.index = le.inverse transform(c matrix.index)
c_matrix.columns = le.inverse_transform(c_matrix.columns)
c matrix
```

9.3.2 Feature Importance

Compare the Random Forest's built-in 'feature importance' with Permutation Feature Importance as documented here.

```
Note
```

🔗 Connections: This next section is the reason you shouldn't blindly run ML algorithms on your da

```
mdi_importances = pd.Series(
  rfc.feature_importances_, index=rfc.feature_names_in_
).sort values(ascending=True)
ax = mdi importances.plot.barh()
ax.set_title("Random Forest Feature Importances (MDI)")
ax.figure.tight_layout()
```

9.3.3 Permutation Feature Importance

```
result = permutation_importance(
  rfc, X_test, y_test, n_repeats=10, random_state=42, n_jobs=2
sorted_importances_idx = result.importances_mean.argsort()
importances = pd.DataFrame(
  result.importances[sorted_importances_idx].T,
  columns=X test.columns[sorted importances idx],
)
ax = importances.plot.box(vert=False, whis=10)
ax.set_title("Permutation Importances (Test Set)")
ax.axvline(x=0, color="k", linestyle="--")
ax.set xlabel("Decrease in accuracy score")
ax.figure.tight_layout()
```

9.4 Shapely Values

Shapely values are a big part of explainable AI and they work (very broadly) by permuting the data to explore how sensitive the predictions made by the model are to the results that you see. For these we need to install two libraries: shap (to do the heavy lifting) and slicer to deal with the data.

9.4.1 Install Libraries

We should now have this already available in Docker, but just in case...

```
try:
import shap
except ModuleNotFoundError:
! pip install slicer shap
import shap
```

9.4.2 Check for Data Types

You are looking for anything other than int64 or float64 for the most part. Boolean should be fine, but pandas' internal, nullable integer type will give you a ufunc error.

```
X_test.info()
X_test['beds'] = X_test.beds.astype('int')
9.4.3 Plot Partial Dependence
shap.partial_dependence_plot(
   "price", rfc.predict, X_test, ice=False,
   model_expected_value=True, feature_expected_value=True
)
```

9.4.4 Calculate Shapely Values

This can take a *long* time: 4-5 hours (!!!) without developing a *strategy* for tackling it. See the long discussion here. I've taken the approach of subsetting the data substantially (the model is already trained so it won't impact the model's predictions) with a 20% fraction of the test data and an explainer sample of 5%. On my laptop the 'Permutation explainer' stage took about 14 minutes, but your results may obviously be rather different.

```
Xsample = shap.utils.sample(X_test.sample(frac=0.2, random_state=41), 10) explainer = shap.Explainer(rfc.predict, Xsample)
```

Now we calculate the shap values for the 5% sample from X_test.

```
6 Caution
```

⚠ Warning: This next block is the one that takes a long time to run. I got between 3mn and 4mn.

```
%%time
shap_values = explainer(X_test.sample(frac=.05, random_state=42))
```

9.4.5 Single Observation

Now you can take any random record (sample_ind) and produce a shap plot to show the role that each attribute played in its classification. Note that getting these plots to save required some searching on GitHub.

```
sample_ind=250
shap.plots.waterfall(shap_values[sample_ind], max_display=14, show=False);
plt.title(f"Shapely values for observation #{sample_ind} ({X_test.sample(frac=.05, random_state=42).iloc[sample_i plt.tight_layout()
#plt.savefig('practical-09-waterfall.png', dpi=150)
```

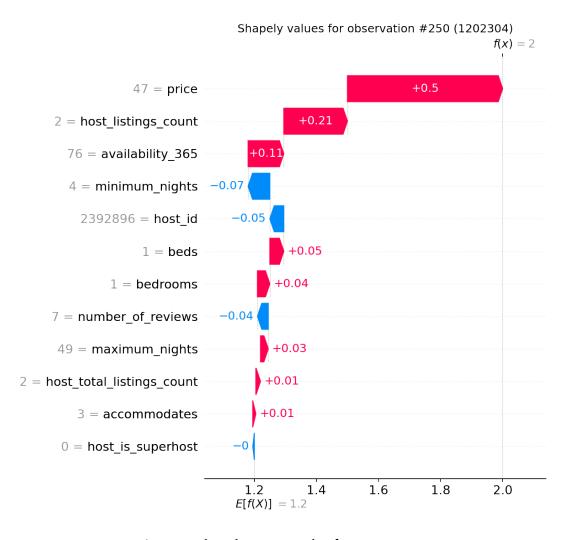


Figure 1: Shapely Feature Plot for Feature 250

9.4.6 All Observations

shap.plots.beeswarm(shap_values, show=False) plt.title(f"Shapely Swarm Plot for Sample")

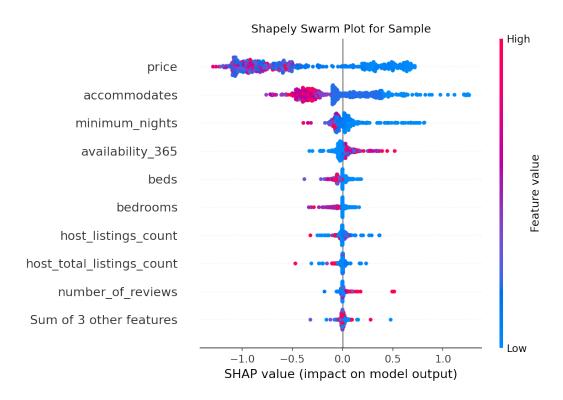


Figure 2: Shapely Swarm Plot

9.5 Wrap-Up

- Find the appropriate eps value: Nearest Neighbour Distance Functions or Interevent Distance Functions
- Clustering Points
- Regionalisation algorithms with Aglomerative Clustering

You've reached the end, you're done...

Er, no. This is barely scratching the surface! I'd suggest that you go back through the above code and do three things: 1. Add a lot more comments to the code to ensure that really have understood what is going on. 2. Try playing with some of the parameters (e.g. my thresholds for skew, or non-normality) and seeing how your results change. 3. Try outputting additional plots that will help you to understand the *quality* of your clustering results (e.g. what *is* the makeup of cluster 1? Or 6? What has it picked up? What names would I give these clsuters?).

If all of that seems like a lot of work then why not learn a bit more about machine learning before calling it a day?

See: Introduction to Machine Learning with Scikit-Learn.