

# Assignment 5 DATA 301

1)

Age	Income	Owng Car	Distance
1	2	Yes	1
3	3	Yes	1.5
1	1	No	1.5
2	3	Yes	1
1	3	No	1.5
3	2	No	1
2	1	Yes	1
1	1	Yes	1.5
3	2	No	1
3	3	Yes	1.5

A person who is middle age, middle class owns a car  
(the rows w/ the dashes are the ones I chose)

2) iteration 1

	age	salary	$C1=(23, 22)$	$C2=(40, 80)$	$C3=(70, 30)$
1	23	22	0	60.44	47.67
2	33	50	29.73	30.8	33.6
3	40	80	60.44	0	58.3
4	11	5	20.8	80.41	64.07
5	70	30	47.67	58.3	0

$C1: [1, 2, 4]$

$C2: [3]$

$C3: [5]$

iteration 2

	age	salary	$C1=(21.3, 25.7)$	$C2=(40, 80)$	$C3=(70, 30)$
1	23	22	3.76	60.44	47.67
2	33	50	26.55	30.8	33.6
3	40	80	57.11	0	58.3
4	11	5	23.58	80.41	64.07
5	70	30	47.89	58.3	0

$C1: [1, 2, 4]$

$C2: [3]$

$C3: [5]$



3)

- Can be difficult to predict the number of clusters
- The initial centroids may have a significant impact on the results. That may cause the solution generated from K-means to be a local optimum.

4) Using a neural network to predict words in speech recognition software

5) With regression the output is a real number whereas with Classification the output is either True/False or a choice out of a list of choices

6.) Supervised learning has both testing and training sets whereas unsupervised learning has no training data.

7.) This is useful because sometimes the data doesn't fit nicely to a  $y = mx + b$  equation. For instance, the data could follow a parabola. Using a polynomial transformation fixes this issue.



8.) `model.fit` takes in 2 parameters  $\rightarrow$  the training data's  $x$ -values and the training data's values. The  $x$  values must be a vector.

```
X = np.array([1, 2, 5, 7, 10])  
y = np.array([2, 4, 6, 8, 15])
```

```
model = LinearRegression(fit_intercept = True)  
model.fit(X[:, np.newaxis], y)
```

9.) `make_pipeline` sequentially applies a list of transformations, or in other words, it pipelines a bunch of steps that modify the input.

```
X = np.array([1, 2, 5, 7, 10])  
y = np.array([2, 4, 6, 8, 15])
```

```
model = make_pipeline(PolynomialFeatures(degree=4),  
LinearRegression())
```

```
model.fit(X[:, np.newaxis], y)
```



10.) `model.predict` uses the model and the testing set's  $x$  values to predict the testing set's  $y$  values. In other words, this allows us to predict the future. This should be called after `model.fit`.

11.) The `add` function allows us to add a layer of nodes to the neural network. The `compile` function configures the learning process (i.e. tells it we need to train on binary classification or multi-class classification problem). Also, we provide a loss function (what we're trying to minimize) and a list of metrics.

12.) `model.evaluate` computes the loss, along with the other metrics passed to the `compile` function of the test data.

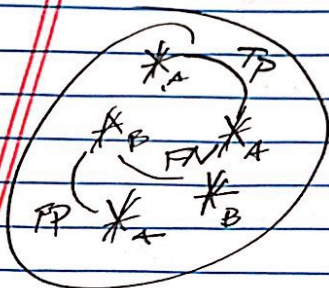
13.) You could use the  $F1$  score.  
The  $F1$  score =  $2 * \frac{\text{precision} * \text{recall}}{\text{precision} + \text{recall}}$ .

$$\text{precision} = \frac{\# \text{ true pos}}{\# \text{ true pos} + \# \text{ false pos}}$$

$$\text{recall} = \frac{\text{true pos}}{\text{true pos} + \text{false neg}}$$

13 (cont) This is a good metric b/c it takes into account each type of possible outcome in its classification.

C.A



14) You could start  $k$  at 0  
and while you don't obtain a  
decent grouping increment  $k$  and run  
the algorithm again.