HW 4: Hierarchical Regression and Bayesian Model Averaging

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On my honor, I have neither given nor received any aid on this assignment.

Part 1: Bayesian Hierarchical Modelling (60 points)

You have been hired by a regional chain of coffee shops to help improve sales and to examine how the personality characteristics of individual store managers might impact the sales numbers of both coffee and food. The client as the following questions they need answered:

- How does conscientiousness and neuroticism impact the sales of coffee and food, and are coffee and food impacted differently?
- Once you control for the personality characteristics of the store managers, what stores should be performing well? (i.e. the rest of the employees might be great, but the store manager might be bringing sales down)
 - 1. Problem Statement What problem are you tackling?
 - 2. Approach Describe the model you are using. Present it both in equation form, as well as a written description of the model. Importantly, you are not writing this for another data scientist, you are writing this for someone who is capable of understanding what a regression is, and what these sorts of models can provide, but has never worked with data analyses or statistics before (so smart, but without the same knowledge base you have.)
 - 3. Prior Rationale List your prior choices and why those were chosen.
 - 4. Findings This is where you present your findings.
 - 5. Summary This is where you summarize and interpret what your analyses uncovered. Again, this is for the client, so it needs to be usable information for them.
 - 6. Diagnostics This is where you put information/plots as to how the estimator performed. This is for technical reference (I like to always have these in the reports I create, but this information is not really for a client per say, more for another data scientist to validate your work.) This doesn't need to be long.

```
In [1]: %matplotlib inline
    import numpy as np
    import pandas as pd
    import matplotlib.pyplot as plt
    import seaborn as sns
    from pymc3 import get_data
```

```
# import as pm to work in Jupyter notebook
import pymc3 as pm
sns.set_context('notebook')
sns.set_style('white')
```

```
In [2]: coffee_sales = pd.read_csv("sales-ds6040sum2021.csv")
    coffee_sales
```

Out[2]:

	sales	food	con	neur	store
0	1.363821	0	0.216553	1.290224	0
1	-1.119747	1	0.216553	1.290224	0
2	-0.180141	0	0.216553	1.290224	0
3	-2.282334	1	0.216553	1.290224	0
4	0.673304	0	0.216553	1.290224	0
•••					
475	0.063563	1	0.306285	-0.070093	19
476	0.206535	0	0.306285	-0.070093	19
477	-1.540875	1	0.306285	-0.070093	19
478	-0.361400	0	0.306285	-0.070093	19
479	-0.937526	1	0.306285	-0.070093	19

480 rows × 5 columns

1.1 - Problem Statement

The problem we are being asked to tackle is to determine for a client (the regional chain of coffee shops) to help improve sales by examinign the personality characteristics of individual store managers. Specifically, we are to determine how conscientiousness and neuroticism impact the sales of coffee and food, and whether coffee and food impacted differently.

1.2 - Approach

As seen from the column headers above, the dataset contains variables related to sales, food/coffee, conscientiousness, neuroticism, and the store identifier. We will need to determine the number of stores and how each specific stores' sales are impacted by the personality of the different store managers.

```
In [3]: print(len(coffee_sales.store.unique()))
20
```

We will be differentiating between 20 different stores, with each individual store appearing to have a consistent conscientiousness and neuroticism level associated witht the store ID number.

Store 14 appears to stand out as a significant leader in sales performance:

```
In [4]: top_10 = coffee_sales.sort_values(by = 'sales', ascending = False).head(10)
top_10
```

```
Out[4]:
```

	sales	food	con	neur	store
349	6.691813	1	0.587351	-0.646036	14
351	5.860888	1	0.587351	-0.646036	14
337	5.413579	1	0.587351	-0.646036	14
353	5.392009	1	0.587351	-0.646036	14
348	5.115312	0	0.587351	-0.646036	14
350	4.934820	0	0.587351	-0.646036	14
345	4.756127	1	0.587351	-0.646036	14
352	4.585693	0	0.587351	-0.646036	14
354	4.516702	0	0.587351	-0.646036	14
343	4.500990	1	0.587351	-0.646036	14

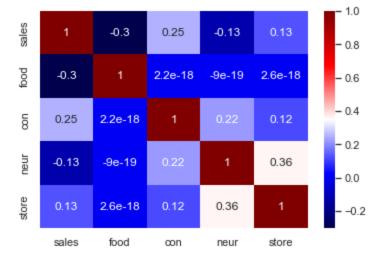
```
In [5]: bottom_10 = coffee_sales.sort_values(by = 'sales').head(10)
bottom_10
```

Out[5]:

	sales	food	con	neur	store
405	-5.651358	1	-0.501073	-0.953021	16
217	-5.384762	1	-0.568532	0.966242	9
301	-4.389198	1	-0.717169	-0.036987	12
183	-4.194727	1	-2.018870	-1.069713	7
227	-3.852984	1	-0.568532	0.966242	9
177	-3.713534	1	-2.018870	-1.069713	7
395	-3.579424	1	-0.501073	-0.953021	16
236	-3.398807	0	-0.568532	0.966242	9
235	-3.365326	1	-0.568532	0.966242	9
239	-3.361530	1	-0.568532	0.966242	9

We also can find what appears to be a positive correlation between conscientiousness and store sales, while store sales and neuroticism appear to have a negative correlation:

```
In [6]: sns.heatmap(coffee_sales.corr(), annot = True, cmap = "seismic")
Out[6]: <AxesSubplot:>
```



For this study, we will be using the hieraarchical modeling approach. In this case, the "hierarchy" is the stores.

Our equation will be:

$$\hat{\text{sales}} = \alpha_{store} + \alpha_f_{store} + \beta * \text{con} + \beta_f * \text{con} + \gamma * \text{neur} + \gamma_f * \text{neur}$$

with:

 β as the effect of conscientiousness for coffee, representing how much the sales are expected to change with a one-unit increase in conscientiousness (con) for coffee products while keeping other variables constant.

 β_f as the difference in the effect of conscientiousness between coffee and food, indicating how much the effect of conscientiousness on sales differs between coffee and food products.

 γ as the effect of neuroticism for coffee. Similar to β , it represents how much the sales are expected to change with a one-unit increase in neuroticism (neur) for coffee products while keeping other variables constant.

 γ_f as the difference in the effect of neuroticism between coffee and food, showing how much the effect of neuroticism on sales differs between coffee and food products.

lpha as the hierarchical effect as well as overall intercept, representing the level of sales at the mean levels of conscientiousness and neuroticism for both coffee and food. Essentially, it gives the baseline sales level when both conscientiousness and neuroticism are at their average values for both coffee and food products.

Code:

```
In [7]: # needed to set these as values to avoid indexing errors

store = coffee_sales.store.values
store_numbers = len(coffee_sales.store.unique())
food = coffee_sales.food.values
con = coffee_sales.con.values
neur = coffee_sales.neur.values
```

In [8]: with pm.Model() as hierarchical_model:

```
# a - overall intercept, level of sales at mean levels of conscientiousness and neur
            mu = pm.Normal('mu a', mu=0., sd=1e5)
            sigma a = pm.HalfCauchy('sigma a', 5)
            # food
            mu a f = pm.Normal('mu a f', mu=0., sd=1e5)
            sigma a f = pm.HalfCauchy('sigma a f', 5)
            # sales intercepts as offsets
            a offset = pm.Normal('a offset', mu=0, sd=1, shape = 20)
            a = pm.Deterministic("a", mu a + a offset * sigma a)
            # food
            a offset f = pm.Normal('a offset f', mu = 0, sd = 1, shape = 20)
            a f = pm.Deterministic("a f", mu a f + a offset f * sigma a f)
            # store level effect of conscientiousness as offset
            b offset = pm.Normal('b offset', mu=0, sd=1)
            b = pm.Deterministic("b", b offset)
            # food
            b offset f = pm.Normal('b offset f', mu=0, sd=1)
            b f = pm.Deterministic("b f", b offset f)
            # store level effect of neuroticism as offset
            c offset = pm.Normal('c offset', mu=0, sd=1)
            c = pm.Deterministic("c", c offset)
            # food
            c offset f = pm.Normal('c offset f', mu=0, sd=1)
            c f = pm.Deterministic("c f", c offset f)
            # Residual Error
            res error = pm.HalfCauchy('res error', 5)
            # This is the specification for the regression equation itself.
            sales reg = a[store] + a f[store]*food + b *con + b f*con*food + c *neur + c f*neur*
            # Data likelihood
            sales likelihood = pm.Normal('sales likelihood', mu = sales reg, sd = res error, obs
In [9]: with hierarchical model:
            hierarchical trace = pm.sample(1000, n init = 50000, tune = 1000)
        /Users/matthewscheffel/opt/anaconda3/lib/python3.9/site-packages/deprecat/classic.py:21
        5: FutureWarning: In v4.0, pm.sample will return an `arviz.InferenceData` object instead
        of a `MultiTrace` by default. You can pass return inferencedata=True or return inference
        data=False to be safe and silence this warning.
         return wrapped (*args , **kwargs )
        Auto-assigning NUTS sampler...
        Initializing NUTS using jitter+adapt diag...
        Multiprocess sampling (4 chains in 4 jobs)
        NUTS: [res error, c offset f, c offset, b offset f, b offset, a offset f, a offset, sigm
        a a f, mu a f, sigma a, mu a]
```

Priors for the fixed effects

- 100.00% [8000/8000 00:10<00:00 Sampling 4 chains, 0

divergences]

```
/Users/matthewscheffel/opt/anaconda3/lib/python3.9/site-packages/scipy/stats/_continuous _distns.py:624: RuntimeWarning: overflow encountered in _beta_ppf return _boost._beta_ppf(q, a, b)
/Users/matthewscheffel/opt/anaconda3/lib/python3.9/site-packages/scipy/stats/_continuous _distns.py:624: RuntimeWarning: overflow encountered in _beta_ppf return _boost._beta_ppf(q, a, b)
/Users/matthewscheffel/opt/anaconda3/lib/python3.9/site-packages/scipy/stats/_continuous _distns.py:624: RuntimeWarning: overflow encountered in _beta_ppf return _boost._beta_ppf(q, a, b)
/Users/matthewscheffel/opt/anaconda3/lib/python3.9/site-packages/scipy/stats/_continuous
```

```
_distns.py:624: RuntimeWarning: overflow encountered in _beta_ppf
  return _boost._beta_ppf(q, a, b)
Sampling 4 chains for 1_000 tune and 1_000 draw iterations (4_000 + 4_000 draws total) t
ook 28 seconds.
```

1.3 - Prior Rationale

The prior distributions for the coefficients α and α_f were chosen as follows:

 μ is set to follow a normal distribution with mean 0 and a large standard deviation of 10000. σ is set to follow a Half-Cauchy distribution with scale parameter 5. An offset is drawn from a normal distribution with mean 0 and standard deviation 1. The shape is 20. The coefficient α is then determined as the sum of μ and each element of the offset vector multiplied by σ . The motivation behind using such wide priors for α and α_f is due to the lack of clear information or strong prior beliefs about their values. By employing these large priors, the model can maintain flexibility and allow for a more general representation of the posterior distributions.

For the coefficients β , β_f , γ , and γ_f , the prior information is as follows:

An offset is drawn from a normal distribution with mean 0 and standard deviation 1. The coefficient of interest (e.g., β , β_f , γ , or γ_f) is determined as the value of the offset. These priors are chosen similarly to the α and α_f priors because there is limited prior information about the data. Moreover, since β , β_f , γ , and γ_f have the same structure within each store (conscientiousness and neuroticism are the same within each store), there is no need for a shape input.

Overall, these prior choices aim to account for uncertainty and avoid imposing strong assumptions on the coefficient values, allowing the model to be more flexible in representing the data's behavior.

1.4 - Findings

43 b_offset_f

45 c_offset_f 0.123

44

88

0.899 0.339

0.269 0.429

0.211

c_offset -0.362 0.278

0.265

-0.867

-0.278

-0.502

```
In [10]: coffee sales summary = pm.summary(hierarchical trace)
         Got error No model on context stack. trying to find log likelihood in translation.
         /Users/matthewscheffel/opt/anaconda3/lib/python3.9/site-packages/arviz/data/io pymc3 3x.
         py:98: FutureWarning: Using `from pymc3` without the model will be deprecated in a futur
         e release. Not using the model will return less accurate and less useful results. Make s
         ure you use the model argument or call from pymc3 within a model context.
          warnings.warn(
In [11]: # grabbing the relevant summary results
         coffee sales summary 1 = coffee sales summary.reset index()
         coffee sales summary 1 = coffee sales summary 1.query("index.str.contains('b|c', case=Fa
         coffee sales summary 1
Out[11]:
                index mean
                               sd hdi_3% hdi_97% mcse_mean mcse_sd ess_bulk ess_tail r_hat
         42
              b_offset 0.269 0.429
                                    -0.502
                                             1.093
                                                         0.010
                                                                 0.007
                                                                         2037.0
                                                                                 2338.0
                                                                                         1.0
```

1.545

0.165

0.523

1.093

0.006

0.006

0.004

0.010

0.005

0.004

0.003

0.007

2822.0

2150.0

3034.0

2037.0

2642.0

2205.0

3021.0

2338.0

1.0

1.0

1.0

```
89
          b_f 0.899 0.339
                                0.265
                                           1.545
                                                        0.006
                                                                   0.005
                                                                            2822.0
                                                                                     2642.0
                                                                                                1.0
90
            c -0.362 0.278
                               -0.867
                                           0.165
                                                        0.006
                                                                   0.004
                                                                            2150.0
                                                                                     2205.0
91
                0.123
                       0.211
                                -0.278
                                           0.523
                                                        0.004
                                                                   0.003
                                                                            3034.0
                                                                                      3021.0
                                                                                                1.0
```

```
In [12]: coffee_sales_summary_2 = coffee_sales_summary.reset_index()
    coffee_sales_summary_2 = coffee_sales_summary_2.query("index.str.contains('14', case=Fal
    coffee_sales_summary_2
```

Out[12]:		index	mean	sd	hdi_3%	hdi_97%	mcse_mean	mcse_sd	ess_bulk	ess_tail	r_hat
	16	a_offset[14]	2.502	0.583	1.456	3.677	0.013	0.009	2115.0	2291.0	1.0
	36	a_offset_f[14]	0.958	0.589	-0.175	2.033	0.010	0.008	3336.0	3212.0	1.0
	62	a[14]	3.585	0.431	2.780	4.407	0.009	0.006	2478.0	2514.0	1.0
	82	a_f[14]	-0.065	0.430	-0.910	0.692	0.007	0.006	3980.0	3319.0	1.0

```
In [13]: coffee_sales_summary_3 = coffee_sales_summary.reset_index()
    coffee_sales_summary_3 = coffee_sales_summary_3.query("index.str.contains('a\[', case=Fa]
    coffee_sales_summary_3.sort_values('mean', ascending=False)
```

Out[13]:		index	mean	sd	hdi_3%	hdi_97%	mcse_mean	mcse_sd	ess_bulk	ess_tail	r_hat
	62	a[14]	3.585	0.431	2.780	4.407	0.009	0.006	2478.0	2514.0	1.0
	60	a[12]	1.699	0.402	1.001	2.530	0.008	0.005	2688.0	2999.0	1.0
	65	a[17]	1.589	0.353	0.982	2.323	0.006	0.005	3105.0	3326.0	1.0
	61	a[13]	1.494	0.499	0.528	2.409	0.010	0.007	2297.0	2723.0	1.0
	52	a[4]	1.250	0.508	0.319	2.224	0.010	0.007	2402.0	2626.0	1.0
	51	a[3]	1.078	0.374	0.354	1.755	0.007	0.005	3066.0	3041.0	1.0
	63	a[15]	1.069	0.289	0.504	1.580	0.004	0.003	4534.0	3149.0	1.0
	55	a[7]	0.646	0.872	-1.055	2.227	0.019	0.014	2093.0	2555.0	1.0
	53	a[5]	0.404	0.443	-0.358	1.280	0.009	0.006	2592.0	2989.0	1.0
	54	a[6]	0.241	0.351	-0.427	0.883	0.006	0.004	3255.0	3493.0	1.0
	66	a[18]	0.201	0.701	-1.074	1.513	0.015	0.011	2229.0	2712.0	1.0
	48	a[0]	0.183	0.436	-0.654	1.009	0.008	0.006	2755.0	3026.0	1.0
	56	a[8]	-0.004	0.349	-0.622	0.687	0.006	0.005	3064.0	2984.0	1.0
	50	a[2]	-0.037	0.404	-0.762	0.764	0.008	0.006	2417.0	2817.0	1.0
	58	a[10]	-0.289	0.324	-0.903	0.307	0.006	0.004	3222.0	3089.0	1.0
	67	a[19]	-0.349	0.292	-0.903	0.192	0.005	0.003	3903.0	3588.0	1.0
	49	a[1]	-0.493	0.645	-1.705	0.687	0.013	0.010	2343.0	2642.0	1.0
	59	a[11]	-0.608	0.345	-1.232	0.058	0.006	0.004	3274.0	3383.0	1.0
	57	a[9]	-0.857	0.474	-1.774	0.018	0.010	0.007	2434.0	2593.0	1.0
	64	a[16]	-1.017	0.406	-1.767	-0.257	0.007	0.005	3095.0	3105.0	1.0

```
In [14]: coffee_sales_summary_4 = coffee_sales_summary.reset_index()
    coffee_sales_summary_4 = coffee_sales_summary_4.query("index.str.contains('a_f', case=Fa
    coffee_sales_summary_4.sort_values('mean', ascending=False)
```

	index	mean	sd	hdi_3%	hdi_97%	mcse_mean	mcse_sd	ess_bulk	ess_tail	r_hat
47	sigma_a_f	0.850	0.192	0.504	1.198	0.004	0.003	2077.0	2584.0	1.0
76	a_f[8]	0.070	0.405	-0.657	0.808	0.006	0.006	5040.0	3228.0	1.0
83	a_f[15]	0.014	0.379	-0.680	0.723	0.005	0.006	4913.0	2676.0	1.0
85	a_f[17]	-0.013	0.401	-0.776	0.732	0.006	0.006	4930.0	3245.0	1.0
81	a_f[13]	-0.031	0.477	-0.952	0.839	0.007	0.006	4664.0	3314.0	1.0
82	a_f[14]	-0.065	0.430	-0.910	0.692	0.007	0.006	3980.0	3319.0	1.0
78	a_f[10]	-0.492	0.394	-1.268	0.215	0.006	0.005	4740.0	2735.0	1.0
74	a_f[6]	-0.557	0.397	-1.280	0.184	0.005	0.005	5223.0	3353.0	1.0
79	a_f[11]	-0.559	0.395	-1.322	0.173	0.005	0.004	5223.0	3166.0	1.0
84	a_f[16]	-0.766	0.426	-1.586	0.033	0.006	0.005	5347.0	2691.0	1.0
1	mu_a_f	-0.864	0.231	-1.300	-0.431	0.005	0.004	2189.0	2068.0	1.0
71	a_f[3]	-0.896	0.416	-1.629	-0.070	0.006	0.004	4867.0	3243.0	1.0
72	a_f[4]	-0.900	0.484	-1.821	-0.005	0.008	0.006	3717.0	3073.0	1.0
69	a_f[1]	-1.046	0.542	-2.024	0.045	0.009	0.007	3674.0	3392.0	1.0
75	a_f[7]	-1.119	0.685	-2.426	0.152	0.013	0.009	2969.0	3054.0	1.0
86	a_f[18]	-1.177	0.569	-2.219	-0.042	0.010	0.007	3384.0	3070.0	1.0
87	a_f[19]	-1.208	0.368	-1.884	-0.489	0.005	0.004	5636.0	3298.0	1.0
77	a_f[9]	-1.265	0.457	-2.199	-0.456	0.007	0.005	4350.0	3348.0	1.0
68	a_f[0]	-1.488	0.431	-2.307	-0.676	0.006	0.005	4401.0	3543.0	1.0
73	a_f[5]	-1.544	0.438	-2.369	-0.712	0.007	0.005	4235.0	3189.0	1.0
70	a_f[2]	-1.944	0.418	-2.708	-1.173	0.006	0.004	4372.0	3142.0	1.0
80	a_f[12]	-2.308	0.452	-3.215	-1.511	0.007	0.005	4099.0	2856.0	1.0

A few takeaways from these findings:

Out[14]:

- Consistent r-hat values around 1 suggest that our sampler performed well, meaning that the sampler has explored the parameter space effectively and provides reliable estimates for the model parameters.
- The means that included food sales tended to outperform the means that did not include food sales. This implies that the introduction of food sales in the stores has contributed positively to the overall revenue of the store.
- As noticed before, store 14 was the best performer in sales, coming in with the highest mean here. However, the store tended to perform well regardless of the impact of the manager's personality.

1.5 - Summary

Impact of Conscientiousness and Neuroticism on Sales of Coffee and Food:

The estimated coefficient for the effect of conscientiousness on coffee sales (β) was approximately 0.27. This means that, on average, a one-unit increase in conscientiousness is associated with a 0.27 increase in coffee sales, keeping other variables constant. The estimated coefficient for the effect of neuroticism on coffee sales (γ) was approximately -0.36. This suggests that, on average, a one-unit increase in neuroticism is associated with a decrease of approximately 0.36 in coffee sales, keeping other variables constant.

The estimated difference in the effect of conscientiousness between coffee and food sales (β_f) was approximately 0.9. This means that the impact of conscientiousness on food sales is stronger than on coffee sales. A one-unit increase in conscientiousness is associated with a higher increase in food sales compared to coffee sales. The estimated difference in the effect of neuroticism between coffee and food sales (γ_f) is approximately 0.12. This suggests that the impact of neuroticism on food sales is weaker than on coffee sales. A one-unit increase in neuroticism is associated with a lower increase in food sales compared to coffee sales.

Overall, conscientiousness and neuroticism have different effects on coffee and food sales. Conscientiousness has a stronger positive impact on food sales than on coffee sales, while neuroticism has a stronger negative impact on coffee sales than on food sales. In general, as seen from the early data visualization and from common sense, a conscientious manager will have a positive impact on sales, while a neurotic manager will have a negative impact on sales.

Stores Performing Well after Controlling for Managers' Personality:

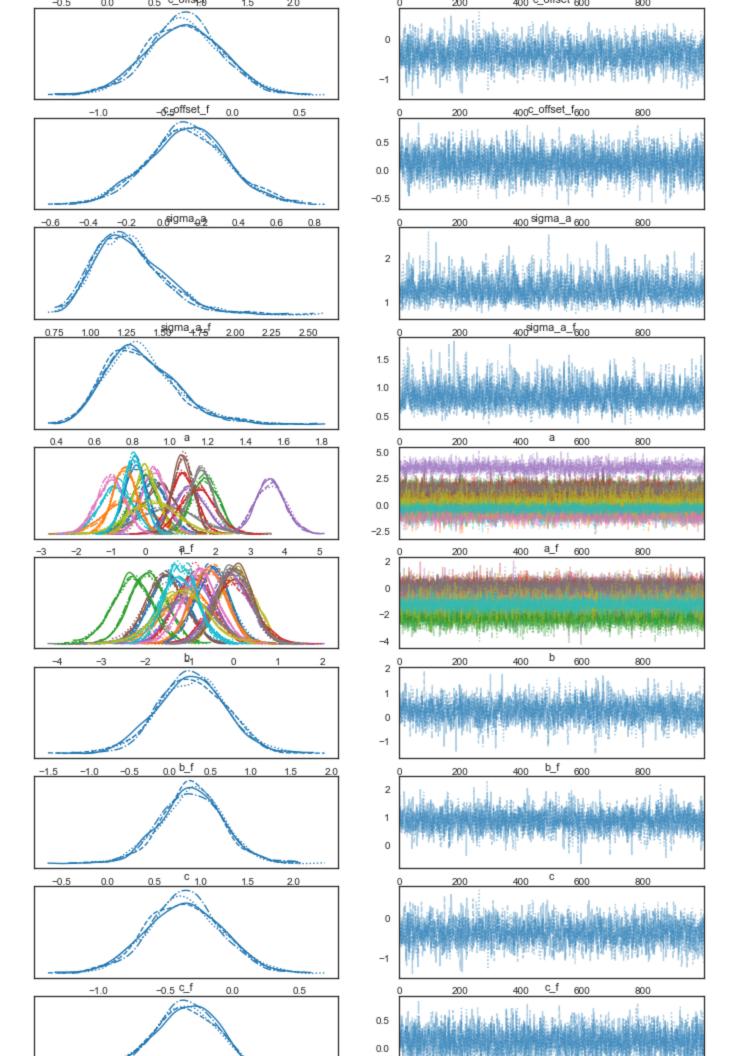
Store 14 stands out as a store that should be performing well once controlling for the manager's personality. Store 14 was the best performer in sales, coming in with the highest mean and tended to perform well regardless of the impact of the manager's personality. Store 14's sales mean was approximately double that of the next closest store, suggesting that it is a high-performing store regardless of the influence of the manager's personality traits. This could indicate that other factors, such as location, customer base, marketing strategies, or product offerings, and other factors play an important role in the store's success, regardless of the impact of the manager's personality on sales performance.

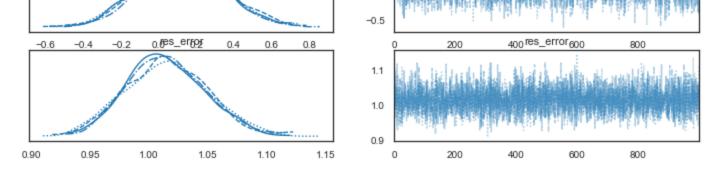
1.6 - Diagnostics

Plots for technical reference:

```
In [15]: pm.plot trace(hierarchical trace)
         Got error No model on context stack. trying to find log likelihood in translation.
         /Users/matthewscheffel/opt/anaconda3/lib/python3.9/site-packages/arviz/data/io pymc3 3x.
         py:98: FutureWarning: Using `from pymc3` without the model will be deprecated in a futur
         e release. Not using the model will return less accurate and less useful results. Make s
         ure you use the model argument or call from pymc3 within a model context.
          warnings.warn(
         Got error No model on context stack. trying to find log likelihood in translation.
         array([[<AxesSubplot:title={'center':'mu a'}>,
Out[15]:
                 <AxesSubplot:title={'center':'mu a'}>],
                [<AxesSubplot:title={'center':'mu a f'}>,
                 <AxesSubplot:title={'center':'mu a f'}>],
                [<AxesSubplot:title={'center':'a offset'}>,
                 <AxesSubplot:title={'center':'a offset'}>],
                [<AxesSubplot:title={'center':'a offset f'}>,
                 <AxesSubplot:title={'center':'a offset f'}>],
                [<AxesSubplot:title={'center':'b offset'}>,
                 <AxesSubplot:title={'center':'b offset'}>],
```

```
[<AxesSubplot:title={'center':'b offset f'}>,
        <AxesSubplot:title={'center':'b offset f'}>],
       [<AxesSubplot:title={'center':'c offset'}>,
        <AxesSubplot:title={'center':'c offset'}>],
       [<AxesSubplot:title={'center':'c offset f'}>,
        <AxesSubplot:title={'center':'c offset f'}>],
       [<AxesSubplot:title={'center':'sigma a'}>,
        <AxesSubplot:title={'center':'sigma a'}>],
       [<AxesSubplot:title={'center':'sigma a f'}>,
        <AxesSubplot:title={'center':'sigma a f'}>],
       [<AxesSubplot:title={'center':'a'}>,
        <AxesSubplot:title={'center':'a'}>],
       [<AxesSubplot:title={'center':'a f'}>,
        <AxesSubplot:title={'center':'a f'}>],
       [<AxesSubplot:title={'center':'b'}>,
       <AxesSubplot:title={'center':'b'}>],
       [<AxesSubplot:title={'center':'b f'}>,
       <AxesSubplot:title={'center':'b f'}>],
       [<AxesSubplot:title={'center':'c'}>,
       <AxesSubplot:title={'center':'c'}>],
       [<AxesSubplot:title={'center':'c f'}>,
        <AxesSubplot:title={'center':'c f'}>],
       [<AxesSubplot:title={'center':'res error'}>,
        <AxesSubplot:title={'center':'res error'}>]], dtype=object)
                                                                          mu_a
                    mu<sub>b_</sub>g_f
                                                                      400 mu_a_f 600
     -0.5
                                      1.5
                                                                                        800
             0.0
                              1.0
                                                             200
                                                   0
                                                  -2
-2.00 -1.75 -1.50 -1.25 -4.00ffset.75 -0.50 -0.25
                                                                      400 a_offset 600
                                                                                        800
                                                  5.0
                                                  25
                                                  0.0
                                                 -2.5
                  o a_offset_f
                                                                      400a_offset_f<sub>600</sub>
                                                  2.5
                                                  0.0
                                                 -2.5
                   1b_offset
                                                                      400 b_offset 600
                                                                                        800
                                                   2
                                                   0
                   b_offset_f_5
                                                                      400b_offset_f600
             -0.5
 -1.5
       -1.0
                               1.0
                                     1.5
                                           2.0
                                                                                        800
```





In [16]: from pymc3 import forestplot, traceplot, plot_posterior
 plt.figure(figsize=(6,14))
 forestplot(hierarchical_trace, var_names=['a'])

ure you use the model argument or call from pymc3 within a model context.

/var/folders/9s/pz872p91741110p_4gz6llzc0000gn/T/ipykernel_7319/3188564304.py:3: Depreca tionWarning: The function `forestplot` from PyMC3 is just an alias for `plot_forest` from ArviZ. Please switch to `pymc3.plot_forest` or `arviz.plot_forest`.

forestplot(hierarchical_trace, var_names=['a'])

Got error No model on context stack. trying to find log_likelihood in translation.

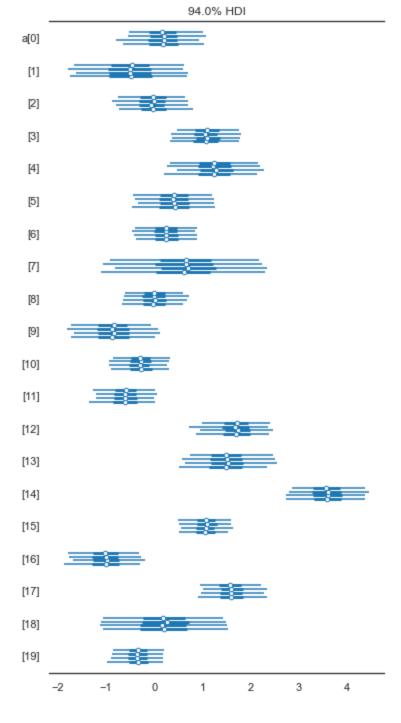
/Users/matthewscheffel/opt/anaconda3/lib/python3.9/site-packages/arviz/data/io_pymc3_3x.

py:98: FutureWarning: Using `from_pymc3` without the model will be deprecated in a futur e release. Not using the model will return less accurate and less useful results. Make s

Out[16]: array([<AxesSubplot:title={'center':'94.0% HDI'}>], dtype=object)

<Figure size 432x1008 with 0 Axes>

warnings.warn(



Part 2: Bayesian Model Averaging with Logistic Regression

- 1. First, revisit your HW2 and calculate the misclassification rate and the cross tabs for 3 variable models that used flat priors that performed best on the testing data. You will have 1 model for LDA and 1 model for QDA.
- 2. Next, use the provided BMA Wine class to fit a Bayesian Model Averaged logistic regression using the training data. Output the variable inclusion probabilities using the summary() function and interpret.
- 3. Finally, obtain the miss-classification rates and cross tabs for the BMA model applied to the training data and the testing data. Compare the performance of the BMA models to the performance of the best LDA and QDA models.

```
In [17]: # load data sets
          wine training = pd.read csv("whitewine-training-ds6040.csv")
          wine testing = pd.read csv("whitewine-testing-ds6040.csv")
In [18]: wine training.head()
                                                                chlorides free.sulfur.dioxide total.sulfur.dioxid
Out[18]:
             fixed.acidity volatile.acidity citric.acid residual.sugar
          0
                0.183032
                             -0.088263
                                         0.223977
                                                       2.798612 -0.038083
                                                                                  0.549032
                                                                                                    0.74478
          1
               -0.640290
                              0.206999
                                        0.056475
                                                      -0.946679
                                                                 0.142355
                                                                                  -1.246502
                                                                                                   -0.14967
          2
                1.476825
                              0.010158
                                        0.558982
                                                      0.092590
                                                                 0.187465
                                                                                  -0.319775
                                                                                                   -0.97351
          3
               -0.757907
                              0.403840
                                       -1.451043
                                                       0.112199 -0.038083
                                                                                  -0.319775
                                                                                                   -0.05552
          4
                0.183032
                             -0.088263
                                        0.223977
                                                       2.798612 -0.038083
                                                                                  0.549032
                                                                                                    0.74478
          wine testing.head()
In [19]:
Out[19]:
             fixed.acidity volatile.acidity citric.acid residual.sugar
                                                                chlorides free.sulfur.dioxide total.sulfur.dioxid
          0
                2.074706
                              2.095780
                                       -0.370398
                                                     -0.442945
                                                               -0.924246
                                                                                 -0.963499
                                                                                                   -0.87840
                              0.031237
                                         0.597571
                0.752463
                                                      -1.000570
                                                                -0.077157
                                                                                  -1.447827
                                                                                                   -0.29067
          2
               -1.050596
                              -0.170181
                                        0.032922
                                                      -1.000570
                                                                 0.016964
                                                                                  -1.750531
                                                                                                   -2.57106
          3
               -1.892024
                               1.541879
                                       -1.822352
                                                      -1.040400 -0.783065
                                                                                  -1.750531
                                                                                                   -2.05386
          4
               -0.569781
                              -0.371600 -0.370398
                                                       1.010863 0.252266
                                                                                 -0.176466
                                                                                                   0.64969
          from mpmath import mp
In [20]:
          import numpy as np
          import pandas as pd
          import statsmodels.api as sm
          from statsmodels.tools import add constant
          import itertools
          from itertools import combinations
          from scipy.stats import multivariate normal
          import matplotlib.pyplot as plt
          import seaborn as sns
          sns.set theme(style = "ticks")
          mp.dps = 50
In [21]: #This class is based on the BMA class provided by Bill Basener in: https://www.kaggle.co
          #It has been modified to allow for multinomial regression (logistic regression for more
          #Specifically, I've hardcoded the model as a 3 category multinomial regression, so this
          class BMA Wine:
              def init (self, y, X, **kwargs):
                   # Setup the basic variables.
                   self.y = y
                   self.X = X
                   self.names = list(X.columns)
                   self.nRows, self.nCols = np.shape(X)
                   self.likelihoods = mp.zeros(self.nCols,1)
                   self.likelihoods all = {}
                   self.coefficients mp = mp.zeros(self.nCols,2)
                   self.coefficients = np.zeros((self.nCols, 2))
                   self.probabilities = np.zeros(self.nCols)
                   # Check the max model size. (Max number of predictor variables to use in a model
                   # This can be used to reduce the runtime but not doing an exhaustive sampling.
```

```
if 'MaxVars' in kwargs.keys():
       self.MaxVars = kwargs['MaxVars']
   else:
        self.MaxVars = self.nCols
    # Prepare the priors if they are provided.
    # The priors are provided for the individual regressor variables.
    # The prior for a model is the product of the priors on the variables in the mod
   if 'Priors' in kwarqs.keys():
       if np.size(kwargs['Priors']) == self.nCols:
           self.Priors = kwargs['Priors']
       else:
           print("WARNING: Provided priors error. Using equal priors instead.")
           print("The priors should be a numpy array of length equal tot he number
            self.Priors = np.ones(self.nCols)
   else:
        self.Priors = np.ones(self.nCols)
   if 'Verbose' in kwargs.keys():
       self.Verbose = kwargs['Verbose']
   else:
       self.Verbose = False
   if 'RegType' in kwargs.keys():
       self.RegType = kwargs['RegType']
        self.RegType = 'LS'
def fit(self):
    # Perform the Bayesian Model Averaging
    # Initialize the sum of the likelihoods for all the models to zero.
    # This will be the 'normalization' denominator in Bayes Theorem.
   likelighood sum = 0
    # To facilitate iterating through all possible models, we start by iterating tho
    # the number of elements in the model.
   \max  likelihood = 0
   for num elements in range(1, self.MaxVars+1):
        if self.Verbose == True:
           print("Computing BMA for models of size: ", num elements)
        # Make a list of all index sets of models of this size.
       Models current = list(combinations(list(range(self.nCols)), num elements))
        # Occam's window - compute the candidate models to use for the next iteratio
        # Models previous: the set of models from the previous iteration that satisf
        # Models next: the set of candidate models for the next iteration
        # Models_current: the set of models from Models_next that can be consturcte
                            to a model from Models previous
        # Iterate through all possible models of the given size.
        for model index set in Models current:
            # Compute the linear regression for this given model.
           model X = self.X.iloc[:,list(model index set)]
           model regr = sm.MNLogit(self.y, model X).fit(disp=0)
            # Compute the likelihood (times the prior) for the model.
           model likelihood = mp.exp(-model regr.bic/2)*np.prod(self.Priors[list(mo
           if self.Verbose == True:
                #print("Model Variables:", model index set, "likelihood=", model likeli
            self.likelihoods all[str(model index set)] = model likelihood
```

```
# Add this likelihood to the running tally of likelihoods.
                         likelighood sum = mp.fadd(likelighood sum, model likelihood)
                          # Add this likelihood (times the priors) to the running tally
                          # of likelihoods for each variable in the model.
                         for idx, i in zip(model index set, range(num elements)):
                              self.likelihoods[idx] = mp.fadd(self.likelihoods[idx], model likelih
                              for j in np.arange(model regr.params.shape[1]):
                                  self.coefficients mp[idx,j] = mp.fadd(self.coefficients mp[idx,j
                         max likelihood = np.max([max likelihood, model likelihood]) # get the new
                 # Divide by the denominator in Bayes theorem to normalize the probabilities
                  # sum to one.
                 self.likelighood sum = likelighood sum
                 for idx in range(self.nCols):
                     self.probabilities[idx] = mp.fdiv(self.likelihoods[idx],likelighood sum, pre
                     for j in range(2):
                         self.coefficients[idx,j] = mp.fdiv(self.coefficients mp[idx,j],likeligho
                  # Return the new BMA object as an output.
                 return self
             def predict MAP(self, true class, data):
                 data = np.asarray(data)
                 result = np.zeros((data.shape[0],3))
                 temp = sm.MNLogit(true class, exog=np.asarray(data))
                 result = temp.predict(params = self.coefficients, exog = np.asarray(data))
                 result = pd.DataFrame(result, columns= ["A", "C", "F"])
                 res MAP = result.idxmax(axis=1)
                 to return = pd.DataFrame(('TrueClass':true class, 'MAP':res MAP))
                 return to return
             def misclass rate(self, true class, data):
                 maps = self.predict MAP(true class, data)
                 maps['Mis class'] = maps['MAP'] == maps['TrueClass']
                 mis class = 1 - maps['Mis class'].mean()
                 return mis class
             def misclass xtabs(self, true class, data):
                 maps = self.predict MAP(true class, data)
                 xtabs = pd.crosstab(maps['MAP'], maps['TrueClass'])
                 return xtabs
             def summary(self):
                  # Return the BMA results as a data frame for easy viewing.
                 df = pd.DataFrame([self.names, list(self.probabilities), list(self.coefficients)
                      ["Variable Name", "Probability", "Avg. Coefficient"]).T
                 return df
In [22]: # LDA Class (code given in example from Professor)
          # used in HW 2 & 3
```

```
# used in HW 2 & 3

class LDA():
    def __init__(self, dataset, class_var, priors = None):
        n_class = len(dataset[class_var].unique())
        if priors is None:
            priors = np.repeat(1/n_class, n_class)
```

```
self.priors = np.asarray(priors)
                 self.means = dataset.groupby(class var).mean()
                 self.sigma = dataset.cov()
                 self.class var = class var
                 self.training data = dataset
             def predict probs(self, data = None):
                 if data is None:
                     data = self.training data
                 data temp = data.drop(self.class var, axis = 1)
                 dens list = []
                 col names = []
                 for ind, row in self.means.iterrows():
                     col names.append(ind)
                     dens list.append(multivariate normal.pdf(data temp, mean = np.asarray(row),
                 dens list = pd.DataFrame(np.transpose(np.vstack(dens list)),columns= col names)
                 dens list = dens list.mul(self.priors, axis=1)
                 dens list = dens list.div(dens list.sum(axis=1), axis=0)
                 dens list['True Class'] = data[self.class var]
                 return dens list
             def predict MAP(self, data = None):
                 if data is None:
                     data = self.training data
                 dens list = self.predict probs(data).drop('True Class', axis = 1)
                 map list = dens list.idxmax(axis = 1)
                 maps = {'MAP Class': map list}
                 maps = pd.DataFrame(maps)
                 maps['True Class'] = data[self.class var]
                 return maps
             def misclass rate(self, data = None):
                 if data is None:
                     data = self.training data
                 maps = self.predict MAP(data = data)
                 maps['Mis class'] = maps['MAP Class'] == maps['True Class']
                 mis class = 1 - maps['Mis class'].mean()
                 return mis class
             def misclass xtabs(self, data = None):
                 if data is None:
                     data = self.training data
                 maps = self.predict MAP(data = data)
                 xtabs = pd.crosstab(maps['MAP Class'], maps['True Class'])
                 return xtabs
             def misclass pairplot(self, data = None):
                 if data is None:
                     data = self.training data
                 maps = self.predict MAP(data = data)
                 temp dat = data.copy(deep = True)
                 temp dat['Mis-Classified'] = maps['MAP Class'] != maps['True Class']
                 plot = sns.pairplot(temp dat, hue="Mis-Classified", height = 1.5, aspect = 1.5)
                 return plot
In [23]: # QDA Class (code given in example from Professor)
```

```
In [23]: # QDA Class (code given in example from Professor)
# from HW 2

class QDA(LDA):
    def __init__(self, dataset, class_var, priors = None):
        n_class = len(dataset[class_var].unique())
        if priors is None:
            priors = np.repeat(1/n_class, n_class)
        self.priors = np.asarray(priors)
        self.means = dataset.groupby(class_var).mean()
```

```
gb = dataset.groupby(class var)
   self.sigma = {x: gb.get group(x).cov() for x in gb.groups}
   self.class var = class var
   self.training data = dataset
def predict probs(self, data = None):
   if data is None:
       data = self.training data
   data temp = data.drop(self.class var, axis = 1)
   dens list = []
   col names = []
   for ind, row in self.means.iterrows():
       col names.append(ind)
       dens list.append(multivariate normal.pdf(data temp, mean = np.asarray(row),
   dens list = pd.DataFrame(np.transpose(np.vstack(dens list)),columns= col names)
   dens list = dens list.mul(self.priors, axis=1)
   dens list = dens list.div(dens list.sum(axis=1), axis=0)
   dens list['True Class'] = data[self.class var]
   return dens list
```

Calculate the misclassification rate and the cross tabs for 3 variable models that used flat priors that performed best on the testing data. You will have 1 model for LDA and 1 model for QDA:

LDA:

```
In [24]: # taken from HW 2
In [25]: # loop over all combinations of 3 features, there will be 165 combinations
          def three features combo (priors):
              features combo = [list(x) for x in combinations(wine training.drop('wine quality', a
             LDA training results = []
             LDA testing results = []
              # iterate combos
             for combination in features combo:
                 current columns = combination + ['wine quality']
                 train three = wine training[current columns]
                 test three = wine testing[current columns]
                 LDA 3 = LDA(train three, 'wine quality', priors = priors)
                 train_misclass = LDA_3.misclass rate()
                 test misclass = LDA 3.misclass rate(test three)
                 LDA training results.append((current columns, train misclass))
                 LDA testing results.append((current columns, test misclass))
              return(LDA training results, LDA testing results)
In [26]: # use volatile.acidity, density, and alcohol as variables since they performed best in H
          # flat priors
          flat priors = [1/3, 1/3, 1/3]
          LDA Flat = LDA(wine testing[["volatile.acidity", "density", "alcohol", "wine quality"]],
In [27]: LDA Flat.misclass rate()
```

```
In [28]: LDA Flat.misclass xtabs()
Out [28]: True Class
                         С
                              F
         MAP Class
                 A 49 307
                             46
                   22 413 158
                    9 280 436
         QDA:
In [29]: # taken from HW 2
In [30]: # loop over all combinations of 3 features, there will be 165 combinations
         def QDA three features combo(priors):
             features\_combo\_2 = [list(x) for x in combinations(wine training.drop('wine quality', features\_combo\_2)]
             QDA training results = []
             QDA testing results = []
              # iterate combos
             for combination in features combo 2:
                 current columns = combination + ['wine quality']
                  train three 2 = wine training[current columns]
                  test three 2 = wine testing[current columns]
                  QDA 3 = QDA(train three 2, 'wine quality', priors=priors)
                  train misclass = QDA 3.misclass rate()
                  test misclass = QDA 3.misclass rate(test three 2)
                  QDA training results.append((current columns, train misclass))
                  QDA testing results.append((current columns, test misclass))
              return(QDA training results, QDA testing results)
In [31]: | # use volatile.acidity, density, and alcohol as variables since they performed best in H
          # flat priors
         flat priors = [1/3, 1/3, 1/3]
         QDA Flat = QDA(wine testing[["volatile.acidity", "free.sulfur.dioxide", "alcohol", "wine
         QDA Flat.misclass_rate()
In [32]:
         0.45813953488372094
Out[32]:
In [33]:
         QDA Flat.misclass xtabs()
Out [33]: True Class
                              F
         MAP Class
                 A 49 305
                             30
                    25 399
                            126
                    6 296 484
```

Out[27]: 0.47790697674418603

Use the provided BMA Wine class to fit a Bayesian Model Averaged logistic regression using the training data. Output the variable inclusion probabilities using the summary() function and interpret.

```
In [34]:
          BMA Log Reg = BMA Wine (wine training ['wine quality'], add constant (wine training.iloc[:,
In [35]:
          BMA Log Reg.fit()
          Computing BMA for models of size:
          Computing BMA for models of size: 5
          Computing BMA for models of size:
          Computing BMA for models of size: 10
          Computing BMA for models of size:
          Computing BMA for models of size:
          < main .BMA Wine at 0x7f9aa252fee0>
Out[35]:
In [36]:
          BMA Log Reg.summary()
Out[36]:
                 Variable Name
                               Probability
                                                                          Avg. Coefficient
           0
                                      1.0
                                                    [3.554805025916327, 2.561719389128104]
                         const
                   fixed.acidity
                                 0.050937
                                              [0.006081316362694214, 0.012836499255148182]
           2
                  volatile.acidity
                                      1.0
                                                 [0.2520051321984379, 0.9405138587589458]
           3
                     citric.acid
                                  0.00038
                                            [1.72439437612421e-05, 1.3338342052968865e-05]
           4
                  residual.sugar
                                      1.0
                                                [-0.4297881784368679, -0.8224412258830641]
           5
                                 0.000432
                                           [5.9312647464358614e-05, 6.510663234553125e-05]
                      chlorides
           6
               free.sulfur.dioxide
                                 0.237846
                                              [-0.06733716611616783, -0.09003553653549728]
              total.sulfur.dioxide
                                 0.002822
                                          [0.0005234846449900891, 0.0008579730874491765]
           8
                                 0.238925
                                                   [0.17396196112899043, 0.29392454867368]
                       density
           9
                           рН
                                 0.035454
                                           [-0.008840094487388952, -0.012972530705770762]
           10
                                 0.127061
                                             [-0.006841192294008149, -0.03058931859351857]
                      sulphates
           11
                        alcohol
                                      1.0
                                                 [-0.8159326146405751, -2.031666402821676]
```

The summary provides us with two important outputs: Probability and the Average Coefficient.

Understanding the probabilities and coefficients aids in identifying important predictors that contribute significantly to the model's performance and can help guide us in further analysis.

The probability represents the likelihood that each predictor (variable) will be included in the final true model. A value of 1.0 indicates that the predictor is highly likely to be included, while smaller values suggest lower chances of inclusion. For instance, variables with a probability of 1.0, such as 'volatile.acidity', 'residual.sugar', and 'alcohol', are very likely to be part of the final true model.

The average coefficient for each predictor corresponds to its estimated effect on the target variable in the final true model. It provides insights into the direction and magnitude of the relationship between each predictor and the target variable. Negative coefficients suggest a negative impact on the target variable, while positive coefficients indicate a positive impact.

As mentioned before, 'volatile.acidity', 'residual.sugar', and 'alcohol' have probabilities of 1.0, implying that they are highly likely to be included in the final true model across BMA, LDA, and QDA. These predictors seem to have substantial explanatory power for the target variable. Volatile Acidity and Alcohol were utilized in all 3 of our models (BMA, LDA, and QDA). Conversely, predictors with lower probabilities might have weaker associations with the target variable or may not be consistently selected across the models.

Obtain the miss-classification rates and cross tabs for the BMA model applied to the training data and the testing data. Compare the performance of the BMA models to the performance of the best LDA and QDA models.

BMA Model:

Training:

Testing:

Best LDA:

```
In [47]:
          # use volatile.acidity, density, alcohol
          Best LDA = LDA(wine training[["volatile.acidity", "density", "alcohol", "wine quality"]],
         Best LDA.misclass rate (wine testing [ "volatile.acidity", "density", "alcohol", "wine qua
In [48]:
         0.4877906976744186
Out[48]:
          Best LDA.misclass xtabs(wine testing[["volatile.acidity", "density", "alcohol", "wine qu
In [49]:
Out[49]:
         True Class
                         C
                              F
          MAP Class
                 A 55 335
                             48
                    17 390
                            156
                       275 436
                     8
```

Best QDA:

```
# use volatile.acidity, free.sulfur.dioxide, alcohol
In [50]:
         Best_QDA = QDA(wine_training[["volatile.acidity", "free.sulfur.dioxide", "alcohol", "wine
         Best QDA.misclass rate(wine testing[["volatile.acidity", "free.sulfur.dioxide", "alcohol
In [51]:
         0.4511627906976744
Out[51]:
         Best QDA.misclass xtabs(wine testing[["volatile.acidity", "free.sulfur.dioxide", "alcoho
In [52]:
         True Class
                         C
                              F
Out[52]:
         MAP Class
                 A 48 299
                             35
                   26 437
                    6 264 459
```

Model Comparison:

Among the three models, the BMA model achieved the lowest misclassification rates on both the training and testing datasets, making it the best-performing model in terms of overall classification accuracy. The LDA and QDA models performed relatively worse, with higher misclassification rates compared to the BMA model.

Bayesian Model Averaging (BMA) Model:

The BMA model achieved the lowest misclassification rates on both the training and testing datasets, indicating its strong classification performance compared to the other models. This suggests that the BMA model's ability to average over different models and predictors contributes to its superior performance in handling complex datasets with multiple potential predictors. The cross-tabulations show that the BMA model achieved relatively balanced predictions across all classes, especially for the 'C' and 'F' classes. This balance in predictions might indicate that the BMA model is generalizing well and is less prone to overfitting, which can be particularly important when dealing with imbalanced datasets or complex class distributions.

Linear Discriminant Analysis (LDA) Model:

The LDA model achieved higher misclassification rates compared to the both the BMA and QDA models, suggesting that it might not be capturing the underlying complexities of the dataset as effectively. LDA's linear nature might limit its ability to handle non-linear relationships between predictors and the target variable. The cross-tabulations show that the LDA model exhibits challenges in classifying the 'A' class correctly, with a higher number of misclassifications. This indicates that the LDA model may struggle with distinguishing the 'A' class from the other classes, possibly due to overlapping feature distributions.

Quadratic Discriminant Analysis (QDA) Model:

The QDA model achieved misclassification rates that are lower than LDA but higher than the BMA model. This suggests that QDA's flexibility in allowing different covariance matrices for each class offers some improvement over LDA but might still be less effective than the BMA model in this specific context. The cross-tabulations reveal challenges in correctly classifying the 'A' class, similar to the LDA model. However, the QDA model seems to perform better than LDA in distinguishing the 'A' and 'C' classes, as indicated by the higher number of correct classifications for class 'A' in comparison.

In summary, the comparison and insights reveal that the Bayesian Model Averaging (BMA) model stands out as the best performer among the three models for the classification task at hand. Its ability to handle multiple potential predictors and model uncertainty, along with its balanced predictions, contribute to its superior classification performance. While LDA and QDA have their strengths, the BMA model emerges as our most promising choice when dealing with complex datasets and multiple predictors. Overall, based on the misclassification rates, the BMA model appears to be the most reliable and robust model for this specific classification problem. However, it's also important to consider other factors, such as interpretability and specific objectives, when choosing the best model.

Extra Credit

LDA and QDA

Both Linear Discriminant Analysis (LDA) and Quadratic Discriminant Analysis (QDA) are classification techniques for multinomial logistic regression.

LDA assumes Gaussian-distributed features with equal covariance matrices across classes, resulting in a linear decision boundary. This assumption simplifies the model and makes it computationally efficient, especially in high-dimensional datasets with limited training data. As such, LDA aims to find a linear

combination of features that optimally separates the classes by maximizing the distance between class means and minimizing within-class variance.

Alternatively, QDA relaxes the assumption of equal covariance matrices and allows each class to have its own covariance matrix. This flexibility allows QDA to capture more complex relationships between features and classes with quadratic decision boundaries. However, the cost of this flexibility is increased computational complexity, especially when dealing with high-dimensional data or a large number of classes. QDA seeks to find a quadratic decision boundary that optimally separates the classes, and in doing so, it can capture more intricate and nonlinear data patterns compared to LDA.

The choice between these methods depends on the underlying data distribution and the complexity of relationships between features and classes, with LDA being more computationally efficient but assuming a simpler decision boundary and QDA allowing greater flexibility.

BIC and Bayes Factor:

The Bayesian Information Criterion (BIC) serves as an approximation to the Bayes Factor, which directly quantifies the strength of evidence for one model over another. The Bayes Factor calculates the likelihood of the data under two competing models and expresses the ratio of those likelihoods. A Bayes Factor greater than 1 indicates evidence in favor of one model, while a Bayes Factor less than 1 suggests evidence for the other model. Essentially, the Bayes Factor weighs the relative support for each model based on how well they explain the observed data.

BIC takes a different approach to model comparison that involves computing the log-likelihood of the data given each model, measuring the goodness of fit for the models. However, the BIC also includes a penalty term based on the number of parameters in each model. This penalty term discourages overly complex models by penalizing those with more parameters and favoring simpler explanations when models are equally effective in fitting the data. The balance between the log-likelihood term (goodness of fit) and the penalty term (model complexity) in the BIC determines the final score assigned to each model. The BIC allows for the comparison of different models and selects the one with the lowest BIC value. A lower BIC value indicates a model that strikes an optimal balance between fit and complexity, making it a suitable candidate for the observed data.

The BIC is computationally efficient and works well for for large datasets due to its approximation nature and the simplicity of it. It is often used in machine learning and statistics to help guide model selection. At times, the Bayes Factor provides a more accurate and direct measure of the evidence supporting one model over another, such as when comparing complex models with many parameters. However, Bayes Factors also tend to have higher computational costs than the BIC, making it less practical for usage in a large-scale analysis.