main

September 29, 2024

1 Midterm 1, Spring 2022: USDA Food Labels

Version 1.2

Solution

Changes:

- 1.1 Clarified instructions for exercise 6
- 1.2 Corrected typos in exercise 2 and exercise 4

This problem builds on your knowledge of nested data structures, string processing, and implementation of mathematical functions. It has 7 exercises, numbered 0 to 6. There are 14 available points. However, to earn 100% the threshold is 11 points. (Therefore, once you hit 11 points, you can stop. There is no extra credit for exceeding this threshold.)

Each exercise builds logically on previous exercises, but you may solve them in any order. That is, if you can't solve an exercise, you can still move on and try the next one. Use this to your advantage, as the exercises are **not** necessarily ordered in terms of difficulty. Higher point values generally indicate more difficult exercises.

Code cells starting with the comment ### Loading results load the result of applying the correct solution as described in the markdown cell above it. These can be used to follow along with the overall analysis, but they are not required to be run to move on. Most of these cells are loading rather large data structures (at least in terms of what's human-readable). As such, the data is only loaded - not printed. You are free to use Python to explore it further.

The point values of individual exercises are as follows:

- Exercise 0: 1 point (This one is a freebie!)
- Exercise 1: 1 point
- Exercise 2: 4 points
- Exercise 3: 1 point
- Exercise 4: 2 points
- Exercise 5: 3 points
- Exercise 6: 2 points

We will be working with some data about food. The foods you see in the grocery store are required to have nutrition labels, which provide information to consumers about which nutrients

are present in a food, the amounts present, and the ingredients in that food (listed in order of content). The USDA maintains all of this information and makes it available to the public on their website - we are using the "Global Branded Foods" data from October 2021 in this notebook.

The file linked on the website may change (hopefully this link will work in the future) - https://fdc.nal.usda.gov/fdc-datasets/FoodData_Central_branded_food_json_2021-10-28.zip.

Note to keep the runtimes shorter, we are working with a *sample* of the source data. This will not affect any of the functionality we will be developing in this notebook.

1.1 Exercise 0 (1 Points):

After downloading and unzipping the file, you will find the data stored in a **JSON** format. You may or may not have seen this type of data format before, so we will take care of reading the data into our Python environment. The result of the code below is that food_lod will load the serialized contents of the json file into Python objects - in this case, a list of dicts.

Run the test cell below to load the data. We are treating this as a "test" cell, so you will get one point for just submitting. How generous!

1.2 Exploring the data

Let's start by taking a look at some of the basic attributes about this data.

Well... food_lod is a list, with 30,000 entries, and each of those are of type dict. Let's take a look at some of the keys in one of the dicts.

```
In [3]: food_lod[0].keys()
Out[3]: dict_keys(['foodClass', 'description', 'foodNutrients', 'foodAttributes', 'modifiedData
```

1.3 Exercise 1 (1 Points):

These look like some promising candidates for extracting information about individual foods. There appear to be some "category" related keys, which may be useful for grouping foods and comparing between groups as well. For further analysis, we want to know if the dicts are all of similar structure to the first one. A good start to analyzing this is determining which keys are common to all of them...

Given an input lod, which is a list of dicts complete the function common_keys(lod) to return a Python set of the keys which are common to all of the dicts in lod.

The demo cell below should display the following output:

The cell below will test your solution for Exercise 1. The testing variables will be available for debugging under the following names in a dictionary format. - input_vars - Input variables for your solution. - original_input_vars - Copy of input variables from prior to running your solution. These *should* be the same as input_vars - otherwise the inputs were modified by your solution. - returned_output_vars - Outputs returned by your solution. - true_output_vars - The expected output. This *should* "match" returned_output_vars based on the question requirements - otherwise, your solution is not returning the correct output.

```
In [7]: ### test_cell_ex1

###

### AUTOGRADER TEST - DO NOT REMOVE

###

from tester_fw.testers import Tester_ex1
```

```
tester = Tester_ex1()
for _ in range(20):
    try:
        tester.run_test(common_keys)
        (input_vars, original_input_vars, returned_output_vars, true_output_vars) = telescept:
        (input_vars, original_input_vars, returned_output_vars, true_output_vars) = telescept:
        (input_vars, original_input_vars, returned_output_vars, true_output_vars) = telescept:
        ###
### AUTOGRADER TEST - DO NOT REMOVE
###
print('Passed! Please submit.')
initializing tester_fw.tester_6040
Passed! Please submit.
```

Even if your solution was incorrect or you skipped this exercise, run this cell to see the expected output of a call to common_keys(food_lod).

```
In [8]: ### Loading results
    import pickle
    import os
    path = './resource/asnlib/publicdata/ex1.pkl'
    if not os.path.exists(path):
        with open(path, 'wb') as file:
            pickle.dump(common_keys(food_lod), file)
    with open(path, 'rb') as file:
        food_keys = pickle.load(file)
```

1.4 Exercise 2 (4 Points):

For our analysis, we are interested in the nutritional content and ingredients contained in each food. Additionally we would like to group foods by the categories given. The keys of interest are 'description', 'ingredients', 'labelNutrients', and 'brandedFoodCategory'.

```
'sugars': {'value': 0.0},
'protein': {'value': 2.0},
'calcium': {'value': 0.0},
'iron': {'value': 0.4},
'potassium': {'value': 319},
'addedSugar': {'value': 0.0},
'calories': {'value': 140}},
'brandedFoodCategory': 'Chips, Pretzels & Snacks'}
```

Define extract_basic_data to meet the following requirements. Given a list of dicts, lod, create a new list of dicts called basic_data. For each dict in lod, there should be a corresponding dict in basic_data with the following key/value pairs: - 'description' - str associated with 'description' in the lod dict. - 'list_of_ingredients' - list of all the ingredients associated with 'ingredients' in the lod dict. See "Notes on ingredients" below. - 'raw_nutrients' - dict mapping the nutrient name (str) to it's amount (float). See "Notes on nutrients" below. - 'category' - str associated with 'brandedFoodCategory' in the lod dict.

Notes on ingredients - For each dict, d in lod the **ingredients** are stored as a str associated with the 'ingredients' key.

- Sometimes there is extra information wrapped in parentheses. We do not want to include this information in our analysis, so any text wrapped in () (and the parentheses themselves) should be left out of further processing. There may be **multiple** sets of parentheses in an ingredients string. You can assume that there are not **nested** parentheses. For example strings of this form **will not** occur 'item, item1 (level 1 (another, level)), item2.' The re module may be helpful here. Note that there can be *anything* in between the parentheses and all of that text should be discarded. For example 'ingredient 1, ingredient 2 (ingredient 2.1, ingredient 2.2, [ingredient 2.2.1, ingredient 2.2.2]), ingredient 3.' should result in just ['ingredient 1', 'ingredient 2', 'ingredient 3'] as it's associated 'list_of_ingredients'. The ingredient string ends in '.', which should also be left out of further processing.
- The individual ingredients are separated by ', '. The ingredients in basic_data[i]['list_of_ingredients'] should not have any leading or trailing whitespace.

Notes on nutrients - For each dict, d, in lod, the nutrients are associated with the 'labelNutrients' key. - d['labelNutrients'] is a dictionary of the form {'protein': {'value': 10}, 'riboflavin': {'value': 2}}, i.e. mapping the nutrient name to a dictionary with one key ('value') which is mapped to the amount of that nutrient present in a particular food.

```
In [10]: ### Define extract_basic_data
    def extract_basic_data(lod):
        ###
        #What are we returning?
        basic_data = []

        #loop through each dict in lod
        for label in lod:
            label_dict = {}
        #Extract the keys
```

```
label_dict['description'] = label['description']
              label_dict['category'] = label['brandedFoodCategory']
          #Ingredients -- turns into a list
              clean_ingr = clean_ingredients(label['ingredients']).split(", ")
              label_dict['list_of_ingredients'] = [i.strip() for i in clean_ingr]
          #Nutrients -- just get the value
              label_dict['raw_nutrients'] = {k:v['value'] for k, v in label['labelNutrients']
          #Append the dicts with the extracted keys to basic_data
              basic_data.append(label_dict)
          return basic_data
      #Clean function for cleaning ingredients string
      def clean_ingredients(ing_str):
          #import re (reg ex)
          import re
          #Remove text within ()
          print(ing_str)
          clean = re.sub("\(.*?\)", "", ing_str)#re.sub(pattern, replacement, target_string
          print(clean)
      #Strip whitespace
          clean = clean.strip()
      #remove the final period
          clean = clean[:-1]
          return clean
          ###
The demo cell below should display the following output:
```

```
[{'description': 'KETTLE COOKED POTATO CHIPS, PINK HIMALAYAN SALT & RED WINE VINEGAR',
 'list_of_ingredients': ['POTATOES',
   'VEGETABLE OIL',
  'MALTODEXTRIN',
   'HIMALAYAN SALT',
   'RED WINE VINEGAR',
   'CITRIC ACID',
   'SUGAR',
  'WHITE DISTILLED VINEGAR',
   'NATURAL FLAVOR'],
  'raw_nutrients': {'fat': 7.0,
   'saturatedFat': 0.501,
  'transFat': 0.0,
  'cholesterol': 0.0,
   'sodium': 140.0,
   'carbohydrates': 17.0,
```

```
'fiber': 1.01,
   'sugars': 0.0,
   'protein': 2.0,
   'calcium': 0.0,
   'iron': 0.4,
   'potassium': 319.0,
   'addedSugar': 0.0,
   'calories': 140.0},
  'category': 'Chips, Pretzels & Snacks'},
 {'description': 'TOMATO BASIL PASTA SAUCE',
  'list_of_ingredients': ['TOMATO PUREE',
   'TOMATOES',
   'SUGAR',
   'SOYBEAN OIL',
   'SALT',
   'DRIED ONIONS',
   'DRIED GARLIC',
   'SPICES',
   'LEMON JUICE CONCENTRATE',
   'ROMANO CHEESE'],
  'raw_nutrients': {'fat': 2.0,
   'sodium': 580.0,
   'carbohydrates': 17.0,
   'fiber': 2.94,
   'sugars': 10.0,
   'protein': 3.0,
   'calcium': 29.4,
   'iron': 0.998,
   'potassium': 750.0,
   'addedSugar': 2.05,
   'calories': 89.6},
  'category': 'Prepared Pasta & Pizza Sauces'}]
In [11]: ### define demo inputs
         keys_of_interest = {'description', 'ingredients', 'labelNutrients', 'brandedFoodCateg'
         demo_lod_ex1 = [{k:v for k, v in d.items() if k in keys_of_interest} for d in food_log
In [12]: ### call demo funtion
         extract_basic_data(demo_lod_ex1)
POTATOES, VEGETABLE OIL (CONTAINS ONE OR MORE OF THE FOLLOWING: CANOLA OIL, SAFFLOWER OIL AND/
POTATOES, VEGETABLE OIL, MALTODEXTRIN, HIMALAYAN SALT, RED WINE VINEGAR, CITRIC ACID, SUGAR, V
TOMATO PUREE (WATER, TOMATO PASTE), TOMATOES, SUGAR, SOYBEAN OIL, SALT, DRIED ONIONS, DRIED GA
TOMATO PUREE, TOMATOES, SUGAR, SOYBEAN OIL, SALT, DRIED ONIONS, DRIED GARLIC, SPICES, LEMON J
```

'category': 'Chips, Pretzels & Snacks',
'list_of_ingredients': ['POTATOES',

Out[12]: [{'description': 'KETTLE COOKED POTATO CHIPS, PINK HIMALAYAN SALT & RED WINE VINEGAR'

```
'VEGETABLE OIL',
  'MALTODEXTRIN',
  'HIMALAYAN SALT',
  'RED WINE VINEGAR',
  'CITRIC ACID',
  'SUGAR',
  'WHITE DISTILLED VINEGAR',
  'NATURAL FLAVOR'],
 'raw_nutrients': {'fat': 7.0,
  'saturatedFat': 0.501,
  'transFat': 0.0,
  'cholesterol': 0.0,
  'sodium': 140,
  'carbohydrates': 17.0,
  'fiber': 1.01,
  'sugars': 0.0,
  'protein': 2.0,
  'calcium': 0.0,
  'iron': 0.4,
  'potassium': 319,
  'addedSugar': 0.0,
  'calories': 140}},
{'description': 'TOMATO BASIL PASTA SAUCE',
 'category': 'Prepared Pasta & Pizza Sauces',
 'list_of_ingredients': ['TOMATO PUREE',
  'TOMATOES',
  'SUGAR',
  'SOYBEAN OIL',
  'SALT',
  'DRIED ONIONS',
  'DRIED GARLIC',
  'SPICES',
  'LEMON JUICE CONCENTRATE',
  'ROMANO CHEESE'],
 'raw nutrients': {'fat': 2.0,
  'sodium': 580,
  'carbohydrates': 17.0,
  'fiber': 2.94,
  'sugars': 10.0,
  'protein': 3.0,
  'calcium': 29.4,
  'iron': 0.998,
  'potassium': 750,
  'addedSugar': 2.05,
  'calories': 89.6}}]
```

The cell below will test your solution for Exercise 2. The testing variables will be available for debugging under the following names in a dictionary format. - input_vars - Input variables

for your solution. - original_input_vars - Copy of input variables from prior to running your solution. These *should* be the same as input_vars - otherwise the inputs were modified by your solution. - returned_output_vars - Outputs returned by your solution. - true_output_vars - The expected output. This *should* "match" returned_output_vars based on the question requirements - otherwise, your solution is not returning the correct output.

```
In [13]: ### test_cell_ex2
         ###
         ### AUTOGRADER TEST - DO NOT REMOVE
         ###
         from tester_fw.testers import Tester_ex2
         tester = Tester_ex2()
         for _ in range(20):
             try:
                 tester.run_test(extract_basic_data)
                 (input_vars, original_input_vars, returned_output_vars, true_output_vars) = te
             except:
                 (input_vars, original_input_vars, returned_output_vars, true_output_vars) = to
                 raise
         ###
         ### AUTOGRADER TEST - DO NOT REMOVE
         print('Passed! Please submit.')
initializing tester_fw.tester_6040
XGUNQYBHYDBFWV, D, AQFAPSO, GPJBNDSWOUNERY.
XGUNQYBHYDBFWV, D, AQFAPSO, GPJBNDSWOUNERY.
OEGWZEZ, PUOIWUASFW, PIUSLDKGFGFO.
OEGWZEZ, PUOIWUASFW, PIUSLDKGFGFO.
I, UIVCRP (ZMNPFTMC), OJLRJBWKPROAHW (MJQHDFHLEDBZPQY, HH).
I, UIVCRP, OJLRJBWKPROAHW.
YSMBQKHHVPFRNVH, DXZQVVY, ELPPYTLNPMD, ZPOZMLOVEDLBSOQ, LSCKBEOGFUJTLGS (LINQIBODNU).
YSMBQKHHVPFRNVH, DXZQVVY, ELPPYTLNPMD, ZPOZMLOVEDLBSOQ, LSCKBEOGFUJTLGS .
RJBIBD, JJSKDQBYBVUPGXS (D, PSZKBCBVSTPP, EGGUTELBOQCV), AGCPTGPXFRDDCO, WBSFNHSHNBUDVPT, PR.
RJBIBD, JJSKDQBYBVUPGXS, AGCPTGPXFRDDCO, WBSFNHSHNBUDVPT, PR.
N, GFBBNX, BJARQC.
N, GFBBNX, BJARQC.
ZWNIBFZZQZAXVON, KRZ, SRNVZRKWG, SINZFKCS, YEDVKNJSP (MOKWK, M).
ZWNIBFZZQZAXVON, KRZ, SRNVZRKWG, SINZFKCS, YEDVKNJSP .
HWHLAJOLWQ, WUHYAKIN, UYVIV (PU, W, VAOX), VYBMVWBNYT, BVTNJLMG.
HWHLAJOLWQ, WUHYAKIN, UYVIV, VYBMVWBNYT, BVTNJLMG.
YWHOAZLOGC, THTPEKLTQCMREO, KTH.
YWHOAZLOGC, THTPEKLTQCMREO, KTH.
XIKABJHRBSA, OYSRCJS, D, BS (UETKU), MLH.
XIKABJHRBSA, OYSRCJS, D, BS , MLH.
```

SMLKTYPV, RHELP, LJ, SRYLRUSTB.

SMLKTYPV, RHELP, LJ, SRYLRUSTB.

HDXTTJIYJHFBQF, HMHB (ZLLUTUCDPXLJGS, L), HKBFWSYPIRN, ADDUFZ, YXO.

HDXTTJIYJHFBQF, HMHB , HKBFWSYPIRN, ADDUFZ, YXO.

SP, VWWDSW, UABYCOAXEHLRVLHY.

SP, VWWDSW, UABYCOAXEHLRVLHY.

CXMAO (NXYDKIP, KGBXZYXNCVYL, SP), BEWDNHOSWBNZEWG, YYQ, YYD, TJPOWOOI.

CXMAO , BEWDNHOSWBNZEWG, YYQ, YYD, TJPOWOOI.

OCKT, QGZNUNUWORYURR, HSGH, LBZ.

OCKT, QGZNUNUWORYURR, HSGH, LBZ.

WQUTDLPCPXXI, AWYWPZPAVRYKGTGV, QFP, EQYPOSAGOTP.

WQUTDLPCPXXI, AWYWPZPAVRYKGTGV, QFP, EQYPOSAGOTP.

XGZBESVOF (XHIU, GRX), ERWQMWNCUBHEYPPX, UF.

XGZBESVOF , ERWQMWNCUBHEYPPX, UF.

RSSHUOEJOP, YRIIMZREMQCQ, XNMNMSCNIYAJBNJ, ML.

RSSHUOEJOP, YRIIMZREMQCQ, XNMNMSCNIYAJBNJ, ML.

LLBHQPGSRI, YCTUTNGA, GOCVRZHOLH, INSUCBJHDS.

LLBHQPGSRI, YCTUTNGA, GOCVRZHOLH, INSUCBJHDS.

BPEZBS, RVGALJBADHRDXPLI, TRSXC, AGD.

BPEZBS, RVGALJBADHRDXPLI, TRSXC, AGD.

TRJF, B, IJP, VXCNOYLS (NRJCRUG, RGRP), EHGC (DPWOEXMIMI, FBE).

TRJF, B, IJP, VXCNOYLS, EHGC.

FOFTOIWF, PNKTHSRJ, AKITQKNW.

FOFTOIWF, PNKTHSRJ, AKITQKNW.

JMQ, RXEAVFPVPNTZGCL, DJPGXS, LNZZEFHCZHEMSB.

JMQ, RXEAVFPVPNTZGCL, DJPGXS, LNZZEFHCZHEMSB.

AVDOON, DVDJNUZSGYAPZCON, NIJCXJYHIRN, JGKWHVSIK, TKSCOLOTJUL.

AVDOON, DVDJNUZSGYAPZCON, NIJCXJYHIRN, JGKWHVSIK, TKSCOLOTJUL.

FFFOXW (CRQOVVJY, X), WAGVPKXQUWZFD, VFCOFXOQKFRFOP, EJ.

FFFOXW, WAGVPKXQUWZFD, VFCOFXOQKFRFOP, EJ.

YMTMTWPSPAF, QGLOHOD, IKMMTDMJRE, NXRZSBXJQWHUTFM.

YMTMTWPSPAF, QGLOHOD, IKMMTDMJRE, NXRZSBXJQWHUTFM.

BZELCVAQ (QYLNBX), KEMQIYJQATKR, XUFJ, FZDRG, DBJHHCPXCGIR.

BZELCVAQ, KEMQIYJQATKR, XUFJ, FZDRG, DBJHHCPXCGIR.

ZGQFZ, WUEC, YRMAY.

ZGQFZ, WUEC, YRMAY.

J, NWNULZMNNWXWOJN, IRKFZEVBRRDTJ, N.

J, NWNULZMNNWXWOJN, IRKFZEVBRRDTJ, N.

FKFDMTEOFPNWRHVU, INHQSW, OY.

FKFDMTEOFPNWRHVU, INHQSW, OY.

NXUFYRHJPKXRMFA, RW, JCIQ, GLPIEESADTWVUG, M.

NXUFYRHJPKXRMFA, RW, JCIQ, GLPIEESADTWVUG, M.

SMLKTYPV, RHELP, LJ, SRYLRUSTB.

SMLKTYPV, RHELP, LJ, SRYLRUSTB.

HDXTTJIYJHFBQF, HMHB (ZLLUTUCDPXLJGS, L), HKBFWSYPIRN, ADDUFZ, YXO.

HDXTTJIYJHFBQF, HMHB, HKBFWSYPIRN, ADDUFZ, YXO.

SP, VWWDSW, UABYCOAXEHLRVLHY.

SP, VWWDSW, UABYCOAXEHLRVLHY.

CXMAO (NXYDKIP, KGBXZYXNCVYL, SP), BEWDNHOSWBNZEWG, YYQ, YYD, TJPOWOOI.

CXMAO , BEWDNHOSWBNZEWG, YYQ, YYD, TJPOWOOI.

OCKT, QGZNUNUWORYURR, HSGH, LBZ.

OCKT, QGZNUNUWORYURR, HSGH, LBZ.

GMKRGVVVNIDPXOUR, MQTFWSDMAV, SLOJCRM, PZ, LLRYMOYEGCR.

GMKRGVVVNIDPXOUR, MQTFWSDMAV, SLOJCRM, PZ, LLRYMOYEGCR.

VZ, ESKPVKQJCDJE, XPK, HVQXNZNJDHBAZUJH.

VZ, ESKPVKQJCDJE, XPK, HVQXNZNJDHBAZUJH.

RDNJUKG, PADQTCO (RQKLFJGOCZUFU), NPQWPWRMWTPM, LBEBDADGORNEKQI, GFFMRCA.

RDNJUKG, PADQTCO, NPQWPWRMWTPM, LBEBDADGORNEKQI, GFFMRCA.

AWKJFVARJBOQPZY, FGEMPRZBEVCUIZ, WMWDIMPSCOI, AGUAOE, MGCF.

AWKJFVARJBOQPZY, FGEMPRZBEVCUIZ, WMWDIMPSCOI, AGUAOE, MGCF.

HDBFZPXLR, OHOYEB (OIMASRKJEI, KTUXPYFVENCSIB, AXVCPFQM), SPDTSFMRS.

HDBFZPXLR, OHOYEB, SPDTSFMRS.

HJZBKLM, KEKQPPZWWIFDXA, IPUQUUDK (PVKUE, PEF).

HJZBKLM, KEKQPPZWWIFDXA, IPUQUUDK .

ENFAJYDVFSGOTEYX, ZYOCZ, VZKDFC, HOZMAAUUAJAU.

ENFAJYDVFSGOTEYX, ZYOCZ, VZKDFC, HOZMAAUUAJAU.

IOLF, RGJZQ, WEUYQNGVI, ULEJVMDQSUWORXUN.

IOLF, RGJZQ, WEUYQNGVI, ULEJVMDQSUWORXUN.

ZQ, MLARI, OMDGYFOKPWTGCE, DEMNCVEXTFSERTL, EJPJCKK (EUAJIJKFKHXKOR).

ZQ, MLARI, OMDGYFOKPWTGCE, DEMNCVEXTFSERTL, EJPJCKK .

EBMQMQNKNAOEIOD (WJGMNJQAI, UROUUPGRBFM), BHEFMODATXAOD, ACYJPKGKLMHOJML, UDNZID.

EBMQMQNKNAOEIOD, BHEFMODATXAOD, ACYJPKGKLMHOJML, UDNZID.

W, KSZ, CHGKU, UDNZALEPUGDWSJ, UGBAIOUWSCJTL.

W, KSZ, CHGKU, UDNZALEPUGDWSJ, UGBAIOUWSCJTL.

YBWREDG, Q, UMDR.

YBWREDG, Q, UMDR.

ADUZFAB, BEDUGBKXDHPEV, EJLZFW, I, UQKVOLOADDQIY.

ADUZFAB, BEDUGBKXDHPEV, EJLZFW, I, UQKVOLOADDQIY.

MEVT (PYHPJTHF), OIMQDTGVFZALKEK, FLYGSRLHFYXXLQ.

MEVT, OIMQDTGVFZALKEK, FLYGSRLHFYXXLQ.

SQYJY, SLLEBYLQSSNWXRJ, JBXKWPIHML, Q.

SQYJY, SLLEBYLQSSNWXRJ, JBXKWPIHML, Q.

V, FUBYGIVIX (ORQRW, MNGEEVKZRXD), ZFEZHWIIXBCDYQH.

V, FUBYGIVIX, ZFEZHWIIXBCDYQH.

FITPIWJETWJMCDWI, MO, DPNYQFLVJCFUODQ, SPWGRNMEMPWSUEO (FWCEZIUZPYU, SYIGDEQMXW).

FITPIWJETWJMCDWI, MO, DPNYQFLVJCFUODQ, SPWGRNMEMPWSUEO .

KA, MIF, LBPVGHKMMQNRDSTL, JFQDSCE, SELLNVU.

KA, MIF, LBPVGHKMMQNRDSTL, JFQDSCE, SELLNVU.

GSU (SLXNEIZ), PMC, FPZJCGLA, IWLPNJTKRQKBB.

GSU , PMC, FPZJCGLA, IWLPNJTKRQKBB.

NSTQMASBQSMJ, UQAKGFXJGS, BHUVVD, WKFTOK.

NSTQMASBQSMJ, UQAKGFXJGS, BHUVVD, WKFTOK.

TKUONDFMRIORJPOU, MVRVQJFPPWDOP, BXKYDL (MSWDAGNTMIEVDHPV, DCM), CMNF.

TKUONDFMRIORJPOU, MVRVQJFPPWDOP, BXKYDL, CMNF.

SFYKEOHGVQSHD, VLLBXFDVV, VRPYUQO.

SFYKEOHGVQSHD, VLLBXFDVV, VRPYUQO.

IJMLDAQRFJHFX, WYTVBAUMYVVMX, A (VZZUECSHNHE).

IJMLDAQRFJHFX, WYTVBAUMYVVMX, A .

P, WQUIFN, ADJYRBKDWVOYSFV.

P, WQUIFN, ADJYRBKDWVOYSFV.

GOMYVJVRHF, JCWTASQJBVR, ANYVXOMKVXZHAIPR.

GOMYVJVRHF, JCWTASQJBVR, ANYVXOMKVXZHAIPR.

YTLYX, PDMJKMNKHNDXCLGZ, QW, IMNZFGXGAWDSBJG (CGTQIFJDAMDBZK).

YTLYX, PDMJKMNKHNDXCLGZ, QW, IMNZFGXGAWDSBJG.

XEBV, WHONOTLWNAAV (LUTYYGICCTWIAX), QZXYAMPEJVVLX.

XEBV, WHONOTLWNAAV, QZXYAMPEJVVLX.

GASKEJEXTSL, JWHGPXGVMGRTC, YRXBPZMRRX, HWNRC.

GASKEJEXTSL, JWHGPXGVMGRTC, YRXBPZMRRX, HWNRC.

IFPAFHUSUD, QTG (VOMBE, XP), LY, NQXSJSAGFSDUG.

IFPAFHUSUD, QTG, LY, NQXSJSAGFSDUG.

IFYJWXYBHQHXIRR, TMZSVWBBTCZF, VAGOLCIQMZBDSE, PMKXXSCYKW, QGPGHOYIRF.

IFYJWXYBHQHXIRR, TMZSVWBBTCZF, VAGOLCIQMZBDSE, PMKXXSCYKW, QGPGHOYIRF.

G, DBMWVC, UYTCWOFHAIYKZ (VRYSGGJFCL, FKPGULEBMXDPPBRA).

G, DBMWVC, UYTCWOFHAIYKZ .

EUXOKOSUUWTVXSOY (BXFBXXE, NZ, OMGAT), VIJHLS, SJSGDAVKFOK.

EUXOKOSUUWTVXSOY, VIJHLS, SJSGDAVKFOK.

MYYXMLCXNCUN, TLPUDVCLTYGYKP, OGNPKNTXREINCW, NYZLCXTGTBISVLF (NQMEXYNPARNHNN, OQXBDW, AHJWAKK' MYYXMLCXNCUN, TLPUDVCLTYGYKP, OGNPKNTXREINCW, NYZLCXTGTBISVLF.

SBKZ, UKUVPSOWQOAZRWNU, EMQOTWERAVPGHLFP, CZO (AYXOYPUPSXI), SKYFIAVK.

SBKZ, UKUVPSOWQOAZRWNU, EMQOTWERAVPGHLFP, CZO, SKYFIAVK.

BXMUSYALHM (PW, A, XDSFO), KHRHCSRSNRS, CPAKJORPEHUOP (V), QK.

BXMUSYALHM , KHRHCSRSNRS, CPAKJORPEHUOP , QK.

EKOAYNW, WQJJPWNSUBGHJEVZ, SSZNHJQG, S.

EKOAYNW, WQJJPWNSUBGHJEVZ, SSZNHJQG, S.

EXPNVLVJZA (DVWIUBGHGVKW, ZYJR), ALWHHUIMZVUR, IPSA.

EXPNVLVJZA, ALWHHUIMZVUR, IPSA.

NIYH, I, JYDVCBO, HKIOYGZIBEEHSBQA, K.

NIYH, I, JYDVCBO, HKIOYGZIBEEHSBQA, K.

UWL, BIJ, HNQAJPPQERVN.

UWL, BIJ, HNQAJPPQERVN.

OSUGKSJQ (LZMML), DWM, TMDRFWXLCEREDWXR.

OSUGKSJQ , DWM, TMDRFWXLCEREDWXR.

PAZPL, UTZTCDXMDODA, DZCGCISCUAIM, YTC, ITSACOAKH.

PAZPL, UTZTCDXMDODA, DZCGCISCUAIM, YTC, ITSACOAKH.

FENLQW, MYZISQMGMZRGGLRD, EHHXDR.

FENLQW, MYZISQMGMZRGGLRD, EHHXDR.

GYAHBUNLEMQRE (VLUHGCNHS, ILKGUOLNMTGUDIO), NPN, APIUR.

GYAHBUNLEMQRE, NPN, APIUR.

TFYRAJSUQROGM, FR, CSGXBOFDYUACB, HTTPOCPDULKQQARI, QBJACCRIQFTDTUJ.

TFYRAJSUQROGM, FR, CSGXBOFDYUACB, HTTPOCPDULKQQARI, QBJACCRIQFTDTUJ.

M (URLKMWQYDSDEDX, YZGCCY), QCWHUAYTMVBNYL (BDBDLLPPXUCPF), SFEPXYRIJJYXV, CSUXZYCMKZKZH.

M , QCWHUAYTMVBNYL , SFEPXYRIJJYXV, CSUXZYCMKZKZH.

MBENHVRYHUH, HAPNI, GHPGOVI.

MBENHVRYHUH, HAPNI, GHPGOVI.

```
WLRNEMUTRTVHJXLT, BZUTADFVRVAW, PGRFAHRAKJTARS, RSYVTATD, ERJX. WLRNEMUTRTVHJXLT, BZUTADFVRVAW, PGRFAHRAKJTARS, RSYVTATD, ERJX. Passed! Please submit.
```

Even if your solution was incorrect or you skipped this exercise, run this cell to see the expected output of a call to extract_basic_data(food_lod).

```
In [14]: ### Loading results
    import pickle
    import os
    path = './resource/asnlib/publicdata/ex2.pkl'
    if not os.path.exists(path):
        with open(path, 'wb') as file:
            pickle.dump(extract_basic_data(food_lod), file)
    with open(path, 'rb') as file:
        basic_data = pickle.load(file)
```

1.5 Exercise 3 (1 Points):

Our analysis requires that the foods have at least 3 listed ingredients and have amounts for nutrients: 'fat', 'protein', 'sodium', and 'carbohydrates'.

Given data, a list of dicts, define filter_basic_data(data) to filter out unwanted records.

- This function should return a **new** list containing the same dicts as data with the following exceptions. - You can assume that each dict in data will have 'list_of_ingredients and 'raw_nutrients' as keys and that the respective values for those keys are of type list and dict.

- Any dict with fewer than 3 items in it's 'list_of_ingredients' should not be included. - Any dict, d, where d['raw_nutrients'] does not have **all** of {'fat', 'protein', 'sodium', 'carbohydrates'} as keys should not be included.

```
In [15]: ### Define filter_basic_data
         def filter_basic_data(data):
             new_data = []
             for recipe in data:
                 to_include = {'ingredients':0,
                               'fat': False,
                                'protein': False,
                                'sodium': False,
                               'carbohydrates': False}
                 for nutrient in recipe['raw_nutrients']:
                     to_include[nutrient] = True
                 for ingredient in recipe['list_of_ingredients']:
                     to_include["ingredients"] += 1
                     if to_include["fat"] and to_include["protein"] and to_include["sodium"] a
                         new_data.append(recipe)
             return new_data
             ###
```

The demo cell below should display the following output:

```
[{'list_of_ingredients': [1, 2, 'this ok', 'milk'],
  'raw_nutrients': {'fat': 22,
   'protein': 'caterpillar',
   'carbohydrates': 5,
   'sodium': 100,
   'awesome sauce': 'this one should be kept'}}]
In [16]: ### define demo inputs
         demo_data_ex3 = [
             {
                 'list_of_ingredients': [1, 2, 'this ok', 'milk'],
                 'raw_nutrients': {
                     'fat':22,
                      'protein': 'caterpillar',
                      'carbohydrates': 5,
                      'sodium': 100,
                      'awesome sauce': 'this one should be kept'}
             },
             {
                 'list_of_ingredients': ['catfish', 2, 'cse6040', 'bicycle'],
                 'raw_nutrients': {
                     'fat':12,
                      'carbohydrates': 35,
                      'sodium': 70,
                      'awesome sauce': 'this one should be rejected - no protein'
                 }
             },
                 'list_of_ingredients': ['marble', 2.5],
                 'raw_nutrients': {
                      'fat':12,
                      'carbohydrates': 35,
                      'protein': 7,
                      'sodium': 70,
                      'awesome sauce': 'this one should be rejected too - not enough ingredient
                 }
             }
         ]
In [17]: ### call demo funtion
         filter_basic_data(demo_data_ex3)
Out[17]: [{'list_of_ingredients': [1, 2, 'this ok', 'milk'],
           'raw_nutrients': {'fat': 22,
            'protein': 'caterpillar',
            'carbohydrates': 5,
```

```
'sodium': 100,
'awesome sauce': 'this one should be kept'}}]
```

The cell below will test your solution for Exercise 3. The testing variables will be available for debugging under the following names in a dictionary format. - input_vars - Input variables for your solution. - original_input_vars - Copy of input variables from prior to running your solution. These *should* be the same as input_vars - otherwise the inputs were modified by your solution. - returned_output_vars - Outputs returned by your solution. - true_output_vars - The expected output. This *should* "match" returned_output_vars based on the question requirements - otherwise, your solution is not returning the correct output.

```
In [18]: ### test_cell_ex3
         ###
         ### AUTOGRADER TEST - DO NOT REMOVE
         ###
         from tester_fw.testers import Tester_ex3
         tester = Tester_ex3()
         for _ in range(20):
             try:
                 tester.run_test(filter_basic_data)
                 (input_vars, original_input_vars, returned_output_vars, true_output_vars) = to
             except:
                 (input_vars, original_input_vars, returned_output_vars, true_output_vars) = to
                 raise
         ###
         ### AUTOGRADER TEST - DO NOT REMOVE
         ###
         print('Passed! Please submit.')
initializing tester_fw.tester_6040
Passed! Please submit.
```

Even if your solution was incorrect or you skipped this exercise, run this cell to see the expected output of a call to filter_basic_data(basic_data).

```
In [19]: ### Loading results
    import pickle
    import os
    path = './resource/asnlib/publicdata/ex3.pkl'
    if not os.path.exists(path):
        with open(path, 'wb') as file:
            pickle.dump(filter_basic_data(basic_data), file)
        with open(path, 'rb') as file:
            filtered_data = pickle.load(file)
```

1.6 Exercise 4 (2 Points):

We want to compute summary statistics on the nutrients present in each food. While you might be able to find formulas for these statistics and implement them yourselves - there is no need to reinvent the wheel. Feel free to use the statistics module, but you will have to import it yourself.

Given a list of dicts, data structured as basic_data, define make_summary(data, key) to generate a dictionary of summary statistics. - We will assume that each dict in data has a 'raw_nutrients' key mapped to a dictionary which maps nutrients to amounts. - Extract the amount of the nutrient given by key for each dict, d in data - we will call these the observations. I.e. data[0]['raw_nutrients'][key] is one observation. - Note: each entry in data counts as an observation and for all dicts d in data, d['raw_nutrients'][key] is not guaranteed to exist. In such cases, we will interpret the observation as a 0. - Compute statistics on the observations. Store the results in a dictionary with the following mapping: - 'mean' - (float) mean of all observations - 'median' - (float) median of all observations - 'stdev' - (float) population standard deviation of all observations - check your stats notes and documentation to make sure you're computing this correctly - 'min' - (float) minimum - 'max' - (float) maximum

```
In [20]: ### Define make summary
         def make_summary(data, key):
             import statistics
             new_list = []
             for d in data:
                 if key in d['raw_nutrients'].keys():
                     temp = d['raw nutrients'][key]
                     new_list.append(temp)
                 else:
                     new_list.append(0)
             xmean = statistics.mean(new_list)
             xmedian = statistics.median(new_list)
             xstdv = statistics.pstdev(new_list)
             xmin = min(new_list)
             xmax = max(new_list)
             new_dict = {}
             new_dict['mean'] = float(xmean)
             new_dict['median'] = float(xmedian)
             new dict['stdev'] = float(xstdv)
             new_dict['min'] = float(xmin)
             new_dict['max'] = float(xmax)
             return new_dict
```

```
###
#d = data
#print(d)
#new_list =
#nutrient_dict = d[0]['raw_nutrients'][key] #, d[1]['raw_nutrients'][key]
#print(nutrient_dict)
###
```

The demo cell below should display the following output:

```
key: foo
{'mean': 18.714285714285715, 'median': 17.0, 'stdev': 12.75835060672349, 'min': 5.0, 'max': 48
key: bar
{'mean': 23.571428571428573, 'median': 33.0, 'stdev': 14.907880397936646, 'min': 0.0, 'max': 33.0, 'stdev': 0.0, 'max': 33.0, 'stdev': 0.0, 'max': 33.0, 'stdev': 0.0, 'max': 0.0, '
key: baz
{'mean': 100.0, 'median': 100.0, 'stdev': 0.0, 'min': 100.0, 'max': 100.0}
In [21]: ### define demo inputs
                        demo_data_ex4 = [
                                   {'raw_nutrients': {'foo': 12,
                                                                                                                          'bar': 33, 'baz': 100}},
                                   {'raw_nutrients': {'foo': 48,
                                                                                                                           'bar': 33, 'baz': 100}},
                                   {'raw_nutrients': {'foo': 17,
                                                                                                                          'bar': 33, 'baz': 100}},
                                   {'raw_nutrients': {'foo': 5,
                                                                                                                                                           'baz': 100}},
                                   {'raw_nutrients': {'foo': 18,
                                                                                                                                                           'baz': 100}},
                                                                                                                          'bar': 33,
                                                                                                                        'bar': 33, 'baz': 100}},
                                   {'raw_nutrients': {'foo': 12,
                                                                                                                                                           'baz': 100}},
                                   {'raw_nutrients': {'foo': 19,
                        demo_keys_ex4 = ['foo', 'bar', 'baz']
In [22]: ### call demo funtion
                        for k in demo_keys_ex4:
                                   print(f'key: {k}')
                                   print(make_summary(demo_data_ex4, k))
                                   print()
key: foo
{'mean': 18.714285714285715, 'median': 17.0, 'stdev': 12.75835060672349, 'min': 5.0, 'max': 48
key: bar
{'mean': 23.571428571428573, 'median': 33.0, 'stdev': 14.907880397936646, 'min': 0.0, 'max': 3
key: baz
{'mean': 100.0, 'median': 100.0, 'stdev': 0.0, 'min': 100.0, 'max': 100.0}
```

The cell below will test your solution for Exercise 4. The testing variables will be available for debugging under the following names in a dictionary format. - input_vars - Input variables for your solution. - original_input_vars - Copy of input variables from prior to running your solution. These *should* be the same as input_vars - otherwise the inputs were modified by your solution. - returned_output_vars - Outputs returned by your solution. - true_output_vars - The expected output. This *should* "match" returned_output_vars based on the question requirements - otherwise, your solution is not returning the correct output.

```
In [23]: ### test_cell_ex4
         ###
         ### AUTOGRADER TEST - DO NOT REMOVE
         ###
         from tester_fw.testers import Tester_ex4
         tester = Tester_ex4()
         for _ in range(20):
             try:
                 tester.run_test(make_summary)
                 (input_vars, original_input_vars, returned_output_vars, true_output_vars) = to
             except:
                 (input_vars, original_input_vars, returned_output_vars, true_output_vars) = to
                 raise
         ###
         ### AUTOGRADER TEST - DO NOT REMOVE
         ###
         print('Passed! Please submit.')
initializing tester_fw.tester_6040
Passed! Please submit.
In [24]: input_vars
Out[24]: {'data': [{'description': 'clrlenpl',
            'raw_nutrients': {'ubb': 17, 'w': 47, 'oooeulpilsgbur': 17}},
           {'description': 'mmefpcucwgulyxqd',
            'raw_nutrients': {'ubb': 47, 'w': 16, 'oooeulpilsgbur': 32}},
           {'description': 'kpvxexqyuyxol',
            'raw_nutrients': {'ubb': 49, 'w': 35, 'oooeulpilsgbur': 4}},
           {'description': 'wmretynrbnfi', 'raw_nutrients': {'ubb': 6, 'w': 27}},
           {'description': 'kswfxzhkg',
            'raw_nutrients': {'ubb': 6, 'w': 5, 'oooeulpilsgbur': 41}},
           {'description': 'bsoslhxgfdpwgtmb', 'raw_nutrients': {'ubb': 31, 'w': 18}},
           {'description': 'e', 'raw_nutrients': {'w': 27, 'oooeulpilsgbur': 9}},
           {'description': 'hgpevnrn',
            'raw_nutrients': {'ubb': 12, 'w': 26, 'oooeulpilsgbur': 33}},
           {'description': 'ynmkgk',
```

```
'raw_nutrients': {'ubb': 48, 'oooeulpilsgbur': 19}},
           {'description': 'dv',
            'raw_nutrients': {'ubb': 38, 'w': 35, 'oooeulpilsgbur': 33}}],
          'key': 'ubb'}
In [25]: returned_output_vars
Out[25]: {'summary': {'mean': 25.4,
           'median': 24.0,
           'stdev': 18.364095403803585,
           'min': 0.0,
           'max': 49.0}}
In [26]: true_output_vars
Out[26]: {'summary': {'mean': 25.4,
           'median': 24.0,
           'stdev': 18.364095403803585,
           'min': 0.0,
           'max': 49.0}}
```

Even if your solution was incorrect or you skipped this exercise, run this cell. You would get the same result as summary_dict if you were to run the code below with a correct implementation of make_summary.

```
keys = ('fat', 'protein', 'carbohydrates', 'sodium')
{key:make_summary(filtered_data, key) for key in keys}

In [27]: ### Loading results
    import pickle
    import os
    path = './resource/asnlib/publicdata/ex4.pkl'
    if not os.path.exists(path):
        with open(path, 'wb') as file:
            pickle.dump({key:make_summary(filtered_data, key) for key in ('fat', 'protein with open(path, 'rb') as file:
            summary_dict = pickle.load(file)
```

1.7 Exercise 5 (3 Points):

We are interested in whether the amount of one particular nutrient in a food has any relationship with the amounts of other nutrients in the food. For this, we will compare the observations of multiple nutrients and compute the correlation between them.

Given data, a list of dicts, and keys, a list of strings, complete the function create_cor_dict(data, key) to find the correlation between each nutrient listed and all of the other nutrients listed. Return the result as a dict which maps each key to a dictionary mapping the other keys to the correlation between the parent key and the child key. For example, if keys=['fat', 'protein', 'carbohydrates'] then the result would look something like this:

```
{'fat': {
                                                 # parent key is 'fat'
                                                 # parent key is 'fat' --> correlation between
    'protein': 0.1854653535334078,
    'carbohydrates': -0.6720362432582452
                                                 # correlation between 'fat' and 'carbohydrates
 'protein': {
                                                 # 'protein'
     'fat': 0.1854653535334078,
                                                 # 'protein' correlation w/ 'fat'
                                                 # 'protein' correlation w/ 'carbohydrates'
     'carbohydrates': -0.3814834566078096
     },
 'carbohydrates': {
                                                 # 'carbohydrates'
                                                 # 'carbohydrates' correlation w/ 'fat'
     'fat': -0.6720362432582452,
     'protein': -0.3814834566078096}
                                                 # 'carbohydrates' correlation w/ 'protein'
```

You can assume that if d is a dict in data, then d will have 'raw_nutrients' as a key which is mapped to a dict which itself maps strings to integers. For example:

Each dictionary in data should be treated as a single observation. You can compute the correlation with the following formulas. - n is the number of observations. - \bar{x} , \bar{y} - Means nutrient x, and nutrient y. - $\bar{x}y = \frac{1}{n}\sum_{i=0}^{n-1} x_i y_i$ - $\sigma_x =$ **population** standard deviation - check your stats notes and documentation to make sure that you are calculating this correctly - Correlation:

$$c = \frac{\bar{xy} - (\bar{x})(\bar{y})}{\sigma_x \sigma_y}$$

```
In [29]: ### Define make_correlations
    def make_correlations(data, keys):
        ###
        def correlation(x, y):
            from statistics import mean, pstdev
            xy = [x_*y_ for x_, y_ in zip(x, y)]
            return (mean(xy) - mean(x)*mean(y)) / pstdev(x) / pstdev(y)

        def key2list(key):
            return [record['raw_nutrients'].get(key, 0.0) for record in data]

return {k1:{k2: correlation(key2list(k1), key2list(k2)) for k2 in keys if k2 != k
            ###
```

The demo cell below should display the following output:

```
{'foo': {'bar': 0.3328398218980465, 'baz': 0.983194888209125},
  'bar': {'foo': 0.33283982189804656, 'baz': 0.31688680340974},
  'baz': {'foo': 0.983194888209125, 'bar': 0.31688680340974007}}
```

```
In [30]: ### define demo inputs
         ### use naming convention demo_varname_ex_* to name demo variables
         demo_data_ex5 = [
             {'raw_nutrients': {'foo': 12,
                                             'bar': 33, 'baz': 100}},
             {'raw_nutrients': {'foo': 48,
                                            'bar': 33, 'baz': 400}},
             {'raw_nutrients': {'foo': 17,
                                            'bar': 33, 'baz': 150}},
             {'raw_nutrients': {'foo': 5,
                                                         'baz': 35}},
             {'raw_nutrients': {'foo': 18,
                                            'bar': 33, 'baz': 200}},
             {'raw_nutrients': {'foo': 12,
                                            'bar': 33, 'baz': 105}},
             {'raw_nutrients': {'foo': 19,
                                                         'baz': 195}},
         demo_keys_ex5 = ['foo', 'bar', 'baz']
In [31]: ### call demo funtion
        make_correlations(demo_data_ex5, demo_keys_ex5)
Out[31]: {'foo': {'bar': 0.3328398218980465, 'baz': 0.983194888209125},
          'bar': {'foo': 0.33283982189804656, 'baz': 0.31688680340974},
          'baz': {'foo': 0.983194888209125, 'bar': 0.31688680340974007}}
```

The cell below will test your solution for Exercise 5. The testing variables will be available for debugging under the following names in a dictionary format. - input_vars - Input variables for your solution. - original_input_vars - Copy of input variables from prior to running your solution. These *should* be the same as input_vars - otherwise the inputs were modified by your solution. - returned_output_vars - Outputs returned by your solution. - true_output_vars - The expected output. This *should* "match" returned_output_vars based on the question requirements - otherwise, your solution is not returning the correct output.

print('Passed! Please submit.')

```
initializing tester_fw.tester_6040 Passed! Please submit.
```

Even if your solution was incorrect or you skipped this exercise, run this cell. You would get the same result as corr_dict if you were to run the code below with a correct implementation of make_correlations.

```
make_correlations(filtered_data, ('fat', 'carbohydrates', 'protein'))
In [33]: ### Loading results
    import pickle
    import os
    path = './resource/asnlib/publicdata/ex5.pkl'
    if not os.path.exists(path):
        with open(path, 'wb') as file:
            pickle.dump(make_correlations(filtered_data, ('fat', 'carbohydrates', 'protein'))
        with open(path, 'rb') as file:
            corr_dict = pickle.load(file)
        corr_dict
Out[33]: {'fat': {'carbohydrates': 0.42945936057100154, 'protein': 0.5300501948181573},
        'carbohydrates': {'fat': 0.42945936057100154, 'protein': 0.4185373785303934},
        'protein': {'fat': 0.5300501948181573, 'carbohydrates': 0.4185373785303934}}
```

1.8 Exercise 6 (2 Points):

We are interested in the most common ingredients listed for foods. Instead of gathering this information on the whole data set, we want it on a category level. A good strategy for drilling down could be useful for generating category level summaries and correlations as well. The function below will transform our basic_data structure (a list of dictionaries) into a dictionary mapping each category to a list of dictionaries which have that category. Each of these lists will have the same structure as basic_data. You may (or may not) find it useful in completing exercise 6.

Complete the function top_ingredients to accomplish the following: - Parameters - data - list of dicts. You can assume that d['list_of_ingredients'] is a list of strings, and d['category'] is a string - for any d in data. Each of these dicts contains data on a single food. - n - int - number of ingredients to list - We will say that an ingredient's "strength" within a category is given by the following:

 x_i = number of times ingredient x has been listed in position i

Strength_x =
$$3x_0 + 2x_1 + x_2$$

- For each unique category (value of d['category']) compute the strength of all ingredients present in that category. - Return a dictionary mapping each category to a list containing the top n ingredients in that category, ranked by strength in descending order. Only include ingredients which have strength greater than 0. - In the instance of ties (two ingredients having the same strength in a category), break the tie by ranking ingredients alphabetically. - If there are fewer than n ingredients - all of the ingredients should be included. There should always be n or fewer ingredients listed for each category.

The demo cell below should display the following output:

```
{'cat0': ['bar', 'foo', 'baz', 'tux'],
 'cat1': ['bax', 'rak', 'foo'],
 'cat2': ['rah']}
In [36]: ### define demo inputs
         demo_data_ex6 = [
             {'category': 'cat0', 'list_of_ingredients':['foo', 'bar', 'baz', 'tux', 'rak']},
             {'category': 'cat0', 'list_of_ingredients':['bar', 'foo', 'baz', 'tux', 'baz']},
             {'category': 'cat0', 'list_of_ingredients':['bar', 'foo', 'tux']},
             {'category': 'cat0', 'list_of_ingredients':['bar', 'baz', 'tux',]},
             {'category': 'cat1', 'list_of_ingredients':['rak', 'foo', 'bax']},
             {'category': 'cat1', 'list_of_ingredients':['rak', 'bax']},
             {'category': 'cat1', 'list_of_ingredients':['bax', 'rak', 'foo']},
             {'category': 'cat1', 'list_of_ingredients':['bax', 'foo', 'rak']},
             {'category': 'cat2', 'list_of_ingredients':['rah']},
             {'category': 'cat2', 'list_of_ingredients':['rah']},
             {'category': 'cat2', 'list_of_ingredients':['rah']},
         ]
In [37]: ### call demo funtion
         top_ingredients(demo_data_ex6, n=5)
Out[37]: {'cat0': ['bar', 'foo', 'baz', 'tux'],
          'cat1': ['bax', 'rak', 'foo'],
          'cat2': ['rah']}
```

The cell below will test your solution for Exercise 6. The testing variables will be available for debugging under the following names in a dictionary format. - input_vars - Input variables for your solution. - original_input_vars - Copy of input variables from prior to running your solution. These *should* be the same as input_vars - otherwise the inputs were modified by your solution. - returned_output_vars - Outputs returned by your solution. - true_output_vars - The expected output. This *should* "match" returned_output_vars based on the question requirements - otherwise, your solution is not returning the correct output.

```
In [38]: ### test_cell_ex6
         ###
         ### AUTOGRADER TEST - DO NOT REMOVE
         ###
         from tester_fw.testers import Tester_ex6
         tester = Tester_ex6()
         for _ in range(20):
             try:
                 tester.run_test(top_ingredients)
                 (input_vars, original_input_vars, returned_output_vars, true_output_vars) = to
             except:
                 (input_vars, original_input_vars, returned_output_vars, true_output_vars) = to
                 raise
         ###
         ### AUTOGRADER TEST - DO NOT REMOVE
         ###
         print('Passed! Please submit.')
initializing tester_fw.tester_6040
Passed! Please submit.
```

Even if your solution was incorrect or you skipped this exercise, run this cell. You would get the same result as leaders if you were to run the code below with a correct implementation of top_ingredients.

```
top_ingredients(filtered_data, 3)
In [39]: ### Loading results
    import pickle
    import os
    path = './resource/asnlib/publicdata/ex6.pkl'
    if not os.path.exists(path):
        with open(path, 'wb') as file:
            pickle.dump(top_ingredients(filtered_data, 3), file)
        with open(path, 'rb') as file:
            leaders = pickle.load(file)
```

Fin. This is the end of the exam. If you haven't already, submit your work.