Exam BIO134: Programming in Biology HS2020

15.01.2021; 10:30-13:00

Question 1

```
text = 'tgovoyd vlwucqk'
numbers = [3, 9, 7, 0, 9, 5, 7, 8, 2, 8, 3, 6, 7, 0, 6]
```

Write a program that creates a new string based on the string *text* and the list *numbers*. The new string should contain all the letters of *text* at those positions where the integers in the list *numbers* are larger than 5, in the order they occur. It should then print the string:

```
good luck
```

The program should be general enough that it would still work according to the same principle, if the list *numbers* and the string *text* would be of different length and/or contain different values. You may assume that the list *numbers* always contains positive integers, and that it always has the same length as the string *text*.

Question 2

```
rna = 'AUGUUCGAA'
```

Write a program that, based on the string *rna*, creates a dictionary, where the keys are the bases A, U, G, C and the values are lists with numbers that indicate the positions where the respective bases occur in the sequence *rna*.

The program should then print the dictionary:

```
{'A': [0, 7, 8], 'U': [1, 3, 4], 'G': [2, 6], 'C': [5]}
```

Note that a dictionary does not have a specified order, so that the same dictionary can be printed in many different ways. It is only important that every key has the correct associated value.

The program should be general enough that it would still work with another sequence of different length. You may assume that every base appears at least once.

```
import numpy as np
np.random.seed(0)
```

Start your code with the lines above. Then, write a function $n_times_to_threshold()$ that takes two positive integer numbers as input: a threshold and a maximum number. Based on these two input variables, the function should sum up random numbers until the sum of all those random numbers is larger than the threshold. The random numbers should range from 1 to (and including) the maximum number, and should be generated within the function by using np.random.randint(1, max number+1).

The function should return the number of times a random number was generated until the sum exceeded the threshold.

Call the function and print the result using 43 as the threshold and 10 as the maximum number by using the following code:

```
k = n\_times\_to\_threshold(43, 10) print('The '+str(k)+'th number has brought the sum above the threshold!')
```

For the given numbers it should then print:

```
The 9th number has brought the sum above the threshold!
```

Question 4

The file 'fMRI_series.txt' contains two numbers on each line, separated by spaces. The first number is the timepoint in seconds and the second number is the activity level in arbitrary units at that given timepoint, measured by fMRI imaging of a mouse brain.

Starting at timepoint 170s, an external stimulus was presented to the mouse every 60 seconds for 5 times, while continuously recording the activity level every 5 seconds.

Write a program that extracts the data from the file and creates a numpy array *data* with the activity levels as floats, starting with the activity level corresponding to timepoint 170s. The array should have 5 rows (one row per round of stimulation), and 12 columns (one column per timepoint in 60 seconds). Calculate and print the mean of the 5 stimulation rounds for each timepoint during the 60 second period using the following code:

```
import numpy as np #somewhere in your program, before:
print(np.mean(data, 0))
```

It should then print the following array:

```
[13622.021218 13924.207528 13901.032028 13897.297232 13978.789918 13937.638388 13888.950176 13874.45905 13711.046962 13664.24686 13557.45596 13617.378826]
```

The code should be general enough that it would still work according to the same principle if 'fMRI_series.txt' contained a dataset with a different number of stimulation rounds. You may assume that activity levels are always recorded every 5 seconds, stimulations are always 60 seconds apart, and that the first stimulation is invariably started at the timepoint 170s.

Here is a list of our solar system's planets (name, diameter [km], distance from the sun [km]) in alphabetical order:

```
planets = [['Earth', 12742, 149598262], ['Jupiter', 139822, 778340821], ['Mars', 6779, 227943824], ['Mercury', 4878, 57909227], ['Neptune', 49244, 4498396441], ['Saturn', 116464, 1426666422], ['Uranus', 50724, 2870658186], ['Venus', 12104, 108209475]]
```

Write a program that sorts the planets according to their distance from the sun, and prints, based on the sorted list, the exact following lines:

Mercury	4878	km in diameter	57909227	km away from the sun
Venus	12104		108209475	
Earth	12742		149598262	
Mars	6779		227943824	
Jupiter	139822		778340821	
Saturn	116464		1426666422	
Uranus	50724		2870658186	
Neptune	49244		4498396441	

Your program should still work if the list *planets* contained information about a subset of the planets only (some entries in *planets* removed). You may assume that the current output format does not need to be adjusted in that case (eg. width of planet name, space, dots would still equal 10 characters).

Please note, that the points you get for this question are largely based on your printed output. In case not all the printed values are correct, or the order of the planets is incorrect, you may still get part of the points (eg. for the formatting).

```
functional_motifs =
['GAGGTAAAC','TCCGTAAGT','AAGGTTGGA','ACAGTCAGT','TAGGTCATT',
'TAGGTACTG','ATGGTAACT','CAGGTATAC','TGTGTGAGT','AAGGTAAGT']
```

'ACTCAGCCCCAGCGGAGGTGAAGGACGTCCTTCCCCAGGAGCCGGTGAGAAGCGCAGTCGGGGGCACGGGGATGAGCTCAGGGGGCCTCTAGAAAGATGTAGCTGGGACCTCGGGAAGCCCTGGCCTCCAGGTAGTCTCAGGAGAGCTACTCAGGGTCGGGCTTGGGGAGAGGAGGAGGAGGGGGGGTGAGGCCAGCAGCA'

Assume you are interested in finding a transcription factor binding site in a given DNA sequence. The transcription factor binding site is not a strictly defined sequence, but there are some preferred characteristics in a 9 base long motif. You know the sequences of 10 exemplary binding sites for this transcription factor (list *functional_motifs*). A profile, also called a position specific scoring matrix, is a motif descriptor that tries to capture the intrinsic variability characteristic of sequence patterns. Such a profile has 4 rows, one for each of the bases A, C, G and T, and as many columns as the sequence motif length. The profile for the given 10 sequences of length 9 is as follows, where each number in the matrix indicates the frequency with which a given nucleotide has been observed at a given position.

	0	1	2	3	4	5	6	7	8
А	0.4	0.6	0.1	0.0	0.0	0.6	0.7	0.2	0.1
С	0.1	0.2	0.1	0.0	0.0	0.2	0.1	0.1	0.2
G	0.1	0.1	0.7	1.0	0.0	0.1	0.1	0.5	0.1
Т	0.4	0.1	0.1	0.0	1.0	0.1	0.1	0.2	0.6

Write a program to create the profile matrix yourself from the 10 given sequences. The data type you use for this matrix (eg. numpy array, dictionary, list) is up to you. The program should print this matrix such that all frequencies are visible, but the precise formatting and labelling of the matrix is not relevant (simply print your variable).

Then, apply the profile matrix to the sequence *query* and find all the potential binding sites. For this, create a score for every possible subsequence of length 9 by adding up the frequencies of its nucleotides.

Print the subsequences with a score larger than 4.4:

```
position 14: GAGGTGAAG, 4.5 position 127: CAGGTAGTC, 4.5
```

Next, generate one hypothetical, ideal motif that, based on the profile matrix, would have the maximum score, and print it (In this particular example there are two possible ideal motifs. It is enough to find one).

```
AAGGTAAGT or TAGGTAAGT
```

Your program should still work with a different motif, a different number of exemplary binding sites and a different query sequence.

Please note, that the points you get for this question are largely based on your printed output and you will be given 0 points if the program prints nothing. In case not all the printed values are correct, you may still get part of the points.

```
aminoacids=['alanine','cysteine','aspartic acid','glutamic acid',
'phenylalanine','glycine','histidine','isoleucine','lysine',
'leucine','methionine','asparagine','proline','glutamine',
'arginine','serine','threonine','valine','tryptophan','tyrosine']
```

The list *aminoacids* contains the 20 standard amino acids. In addition to their full name, there is a standard one letter code, that uniquely identifies the amino acids.

Write a program to create your own one letter code. Use the following rules for this:

- 1. If the first letter of the name is unique, use it as a one letter code (eg P for Proline)
- 2. For the other amino acids that share the first letter:
 - a. Assign the first letter as code to the amino acid with the shortest name in the group of amino acids with the same first letter (eg. ['Glycine', Glutamic acid', 'Glutamine']: G for Glycine). You may assume that there is exactly one shortest name per group.
 - b. For the remaining amino acids: assign them in alphabetical order to the next free letter in the alphabet (eg. B for Arginine).

The program should print your one letter code in the following manner (note the alphabetical order in the one letter code and the capital letter in the full names):

```
A Alanine
B Arginine
C Cysteine
D Asparagine
E Aspartic acid
F Glutamic acid
G Glycine
H Histidine
I Isoleucine
J Glutamine
K Leucine
L Lysine
M Methionine
N Phenylalanine
O Threonine
P Proline
Q Tryptophan
R ---
S Serine
T Tyrosine
U ---
V Valine
W ---
X ---
Y ---
```

Your program should still work if amino acids were named differently or if there was a different number of amino acids. You may assume that there are not more than 26 amino acids

Please note, that the points you get for this question are largely based on your printed output and you will be given 0 points if the program prints nothing. In case not all the printed values are correct, you may still get part of the points.