session3 clustering solution

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1 [MIRACUM 2019][Session 3] Solution

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- This is prepared for a turorial Data analysis tools (Datenanalysewerkzeuge)

2 Depict a phylogenetic tree of five species

2.0.1 Protocol:

1. Import required libraries 2. Load protein of Hemoglobin subunit alpha data (download) 3. Sequence alignment - calculate the similarities of sequence 4. Visualize the result in dendrogram (phylogenetic tree)

```
[11]: from Bio import SeqIO, Phylo from Bio.Align.Applications import ClustalwCommandline
```

2.0.2 Protocol:

1. Import required libraries 2. Load protein of Hemoglobin subunit alpha data (download) 3. Sequence alignment - calculate the similarities of sequence 4. Visualize the result in dendrogram (phylogenetic tree)

```
[6]: FILE_PATH = 'data/protein.fasta'
records = list(SeqIO.parse(FILE_PATH, "fasta"))
records
```

[6]: [SeqRecord(seq=Seq('MVLSADDKTNIKNCWGKIGGHGGEYGEEALQRMFAAFPTTKTYFSHIDVSPGSA...KYR
 ', SingleLetterAlphabet()), id='rat', name='rat', description='rat Species_rat',
 dbxrefs=[]),

SeqRecord(seq=Seq('MVLSPADKTNVKAAWGKVGAHAGEYGAEALERMFLSFPTTKTYFPHFDLSHGSA...KYR
', SingleLetterAlphabet()), id='human', name='human', description='human
Species_human', dbxrefs=[]),

SeqRecord(seq=Seq('MSLSDTDKAVVKAIWAKISPKADEIGAEALARMLTVYPQTKTYFSHWADLSPGS...KYR
', SingleLetterAlphabet()), id='zebrafish', name='zebrafish',

```
description='zebrafish Species_zebrafish', dbxrefs=[]),
   SeqRecord(seq=Seq('MVLSAADKSNVKAAWGKVGGNAGAYGAEALERMFLSFPTTKTYFPHFDLSHGSA...KYR
', SingleLetterAlphabet()), id='sheep', name='sheep', description='sheep
Species_sheep', dbxrefs=[]),
   SeqRecord(seq=Seq('MVLSPADKTNVKAAWGKVGAHAGEYGAEALERMFLSFPTTKTYFPHFDLSHGSA...KYR
', SingleLetterAlphabet()), id='chimpanzee', name='chimpanzee',
description='chimpanzee Species_chimpanzee', dbxrefs=[])]
```

2.0.3 Protocol:

1. Import required libraries 2. Load protein of Hemoglobin subunit alpha data (download) 3. Sequence alignment - calculate the similarities of sequence 4. Visualize the result in dendrogram (phylogenetic tree)

```
[7]: clustalw_cline = ClustalwCommandline("clustalw2", infile=FILE_PATH)
stdout, stderr = clustalw_cline()
print(stdout)
```

CLUSTAL 2.1 Multiple Sequence Alignments

Sequence format is Pearson

```
Sequence 1: rat
                         142 aa
Sequence 2: human
                         142 aa
Sequence 3: zebrafish
                         143 aa
Sequence 4: sheep
                         142 aa
Sequence 5: chimpanzee
                         142 aa
Start of Pairwise alignments
Aligning...
Sequences (1:2) Aligned. Score:
                                 78
Sequences (1:3) Aligned. Score:
                                 51
Sequences (1:4) Aligned. Score:
                                 76
Sequences (1:5) Aligned. Score:
                                 78
Sequences (2:3) Aligned. Score:
                                 53
Sequences (2:4) Aligned. Score:
                                 86
Sequences (2:5) Aligned. Score:
                                 100
Sequences (3:4) Aligned. Score:
                                 53
Sequences (3:5) Aligned. Score:
                                 53
Sequences (4:5) Aligned. Score:
Guide tree file created:
                           [data/protein.dnd]
There are 4 groups
Start of Multiple Alignment
```

```
Aligning...
```

Group 1: Sequences: 2 Score:3061
Group 2: Sequences: 3 Score:2877
Group 3: Sequences: 4 Score:2764
Group 4: Sequences: 5 Score:2355
Alignment Score 6225

CLUSTAL-Alignment file created [data/protein.aln]

2.0.4 Protocol:

1. Import required libraries 2. Load protein of Hemoglobin subunit alpha data (download) 3. Sequence alignment - calculate the similarities of sequence 4. Visualize the result in dendrogram (phylogenetic tree)

```
[12]: newick_path = 'data/protein.dnd'
tree = Phylo.read(newick_path, "newick")
Phylo.draw_ascii(tree)
```

