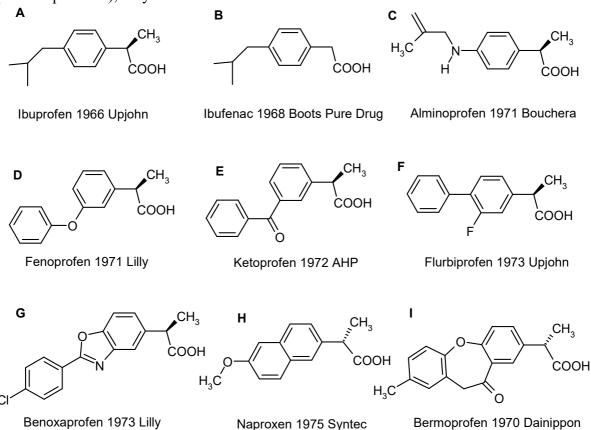
## 5th Assignment

## your name:

1. The analgesic and anti-inflammatory Ibuprofen faced a series of "me-too" drugs shortly after its commercial launch, since the corresponding patent claimed the structure of ibuprofen (see compound **A**), only.



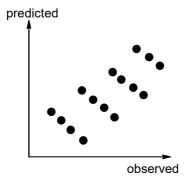
Report the matching compounds(s) (alphabetic character is sufficient) for the given SMARTS 1 to 4. Not all compounds may me matched (5 points for each correct assignment, deduction if more than one is falsely assigned)

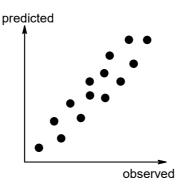
- 1: OC(=O)C(C)c1ccc([C,N])cc1
- 2: OC(=O)C(C)c1cccc([C,O,F])c1
- 3: OC(=O)C(C)c1cccc([#6])c1
- 4: OC(=O)C(C)c1cccc([C,O,#7])c1

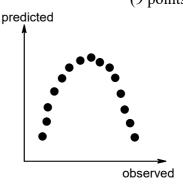
Which compound is not matched by any of the SMARTS strings? (4 points)

2. Mark the correct value of the linear correlation for each of the plots, respectively

(9 points)







$$r^2 = 0.00$$
  
 $r^2 = 0.60$   
 $r^2 = 0.90$   
 $r^2 = 0.95$ 

$$r = -0.98$$
  
 $r = 0.00$   
 $r = 0.90$   
 $r = 0.98$ 

$$r^{2} = 0.00$$
  
 $r^{2} = 0.50$   
 $r^{2} = 0.60$   
 $r^{2} = 0.85$ 

3. Which of the QSAR equations (A or B, respectively C or D) is "better"? (According to which criteria?) (10 points)

*n*: number of molecules;  $r^2$ : squared correlation coefficient;  $q^2$ : cross-validated  $r^2$ ; s: standard error

- A)  $\log 1/K_i = 0.26(\pm 0.14) \text{ClogP} 1.36(\pm 0.43) \text{NHB} + 9.66(\pm 0.39) \text{HL}$  where n = 21,  $r^2 = 0.853$ ,  $q^2 = 0.746$ , s = 0.326
- B)  $\log 1/K_i = 0.34(\pm 0.13)\text{ClogP} 0.96(\pm 0.27)\text{MR} + 0.85(\pm 0.19)\text{B5Y}$ where n = 21,  $r^2 = 0.853$ ,  $q^2 = 0.807$ , s = 0.302
- C) log 1/C =  $0.26(\pm 0.18)$ ClogP  $1.36(\pm 0.43)$ MR  $+ 9.66(\pm 0.39)$ B5Y where n = 16,  $r^2 = 0.821$ ,  $q^2 = 0.794$ , s = 0.354
- D) log 1/C =  $0.23(\pm 0.15)$ ClogP  $1.20(\pm 0.56)$ MR  $+ 10.74(\pm 0.67)$ B5Y  $0.62(\pm 0.12)$ NN4  $+ 0.14(\pm 0.05)$ GEOM6 where n = 16,  $r^2 = 0.892$ ,  $q^2 = 0.632$ , s = 0.325
- 4. Chemical alterations shall be carried for the shown compounds. Draw a new structure that includes the suitable modifications. (12 points)

- a) suggest a modification that reduces the number of freely rotatable bonds
- b) suggest a modification that will lead to increased metabolic stability of the lower benzene ring