

Corrigé TD n° 8 : Structures des molécules organiques

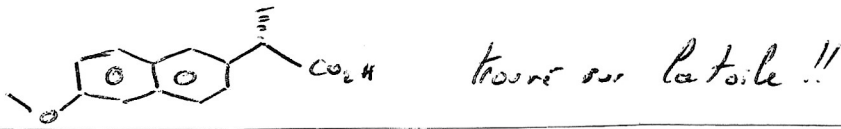
Exercice 1: $C_x H_y O_z$

$$\%C = \frac{12,0x}{230,3} \Rightarrow x = 14$$

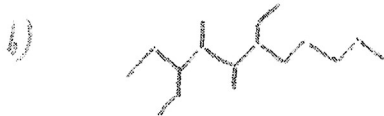
$$\%H = \frac{1,01y}{230,3} \Rightarrow y = 14$$

$$\%O = 1 - \%C - \%H = \frac{16z}{230,3} \Rightarrow z = 3$$

Naproxen: $C_{14}H_{14}O_3$



Ex 2: a) 4-ethyl-2,2,6,7-tetraméthyldecane

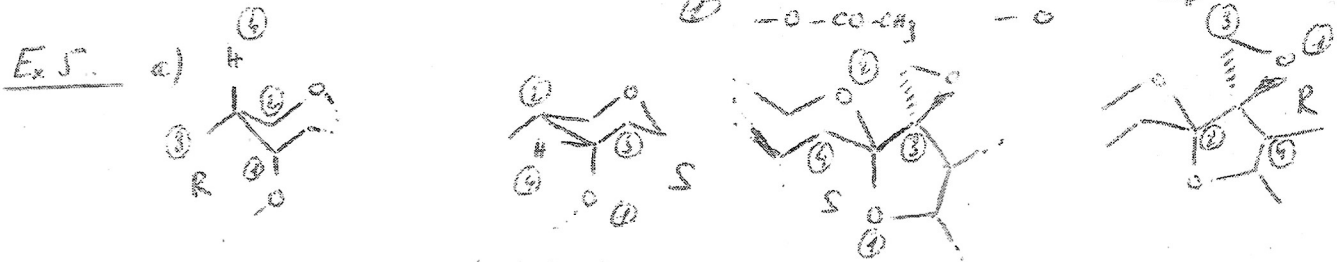


Ex 3 $DI = \frac{4+1-16}{2} = 0$

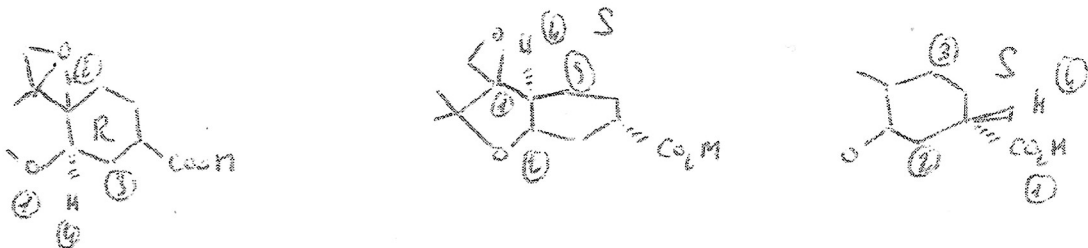


Ex 4

	1 ^{er} rang	2 ^e rang	3 ^e rang	4 ^e rang	5 ^e rang
① -OCH ₃	-O	-C	-C	-C	-C
② -CH	-O	-H	-H	-H	-H
③ -CH ₂ -O-CH ₃	-C	-C	-C	-C	-C



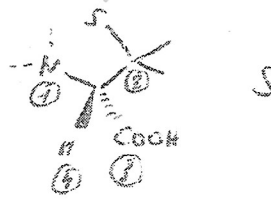
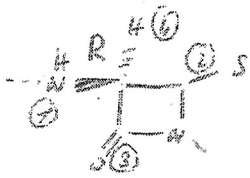
Il y a 7 carbones asymétriques de la phénylthiocarbamide.



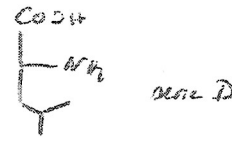
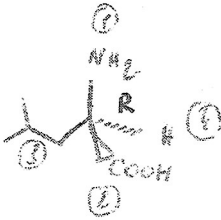
7 C* et pas de contraintes particulières donc $2^7 = 128$ stéréoisomères de configuration possible

b) 11/2 a 3 carbonos asintéticos.

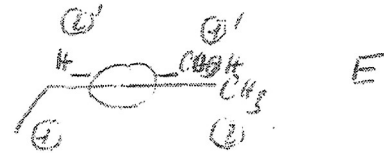
para de construcción para la construcción $2^3 = 8$
 iteraciones de configuración posible



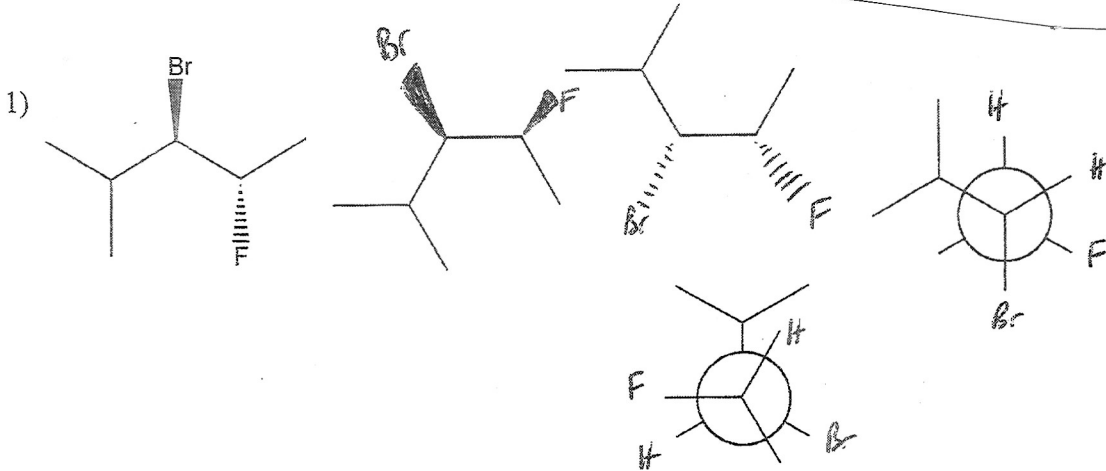
Ex 6



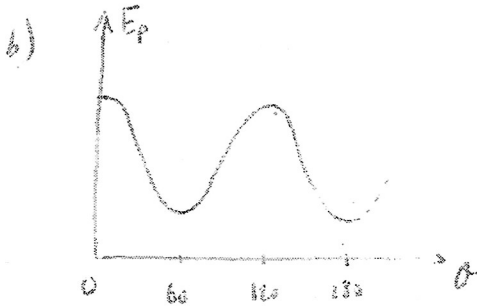
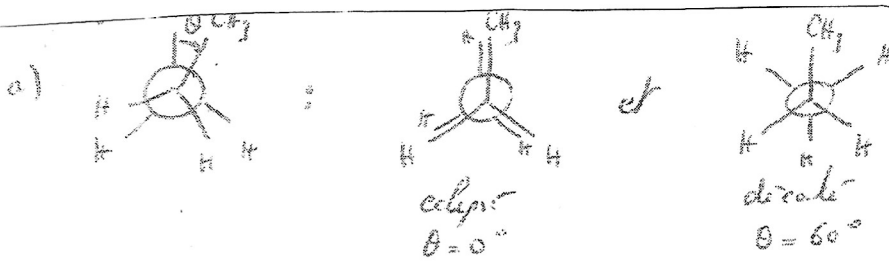
Ex 7



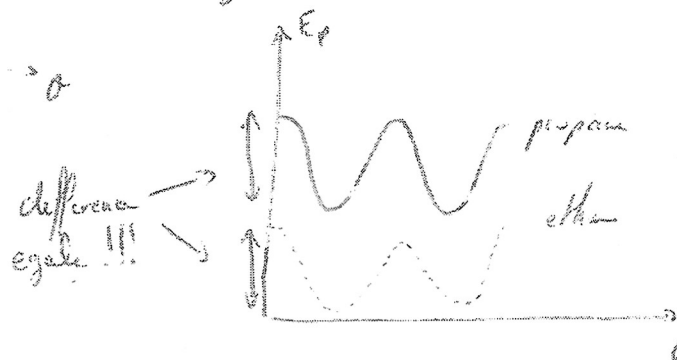
Exercice 8:



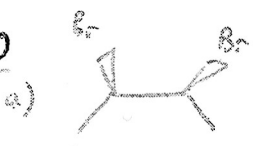
Exercice 9:



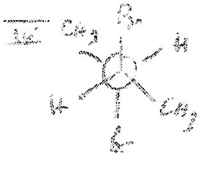
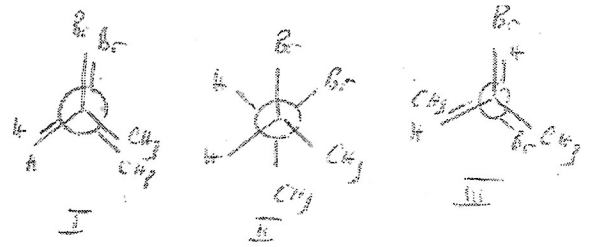
c) EVIDEMMENT, l'interaction gauche H/CH₃ est + désfavorable que l'interaction gauche H/H mais la traverse.



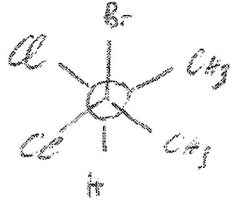
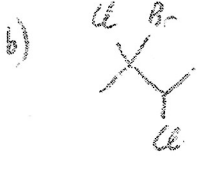
Ex 10



Conformation particulière:

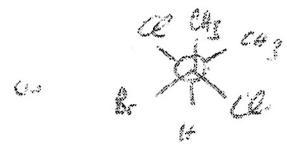
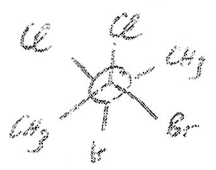


Il s'agit du conformation la + stable: diéclipte et peut favoriser le moment dipolaire associé aux liaisons C-Br très polaires

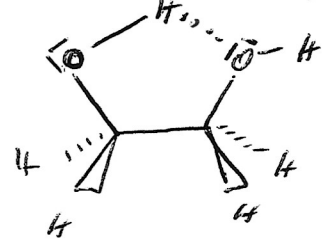


La conformation la + stable correspond à un conformation diéclipte où les moments dipolaires sont en part ant.

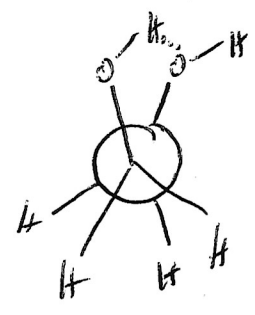
Il y a 2 possibilités



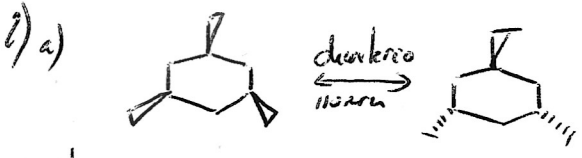
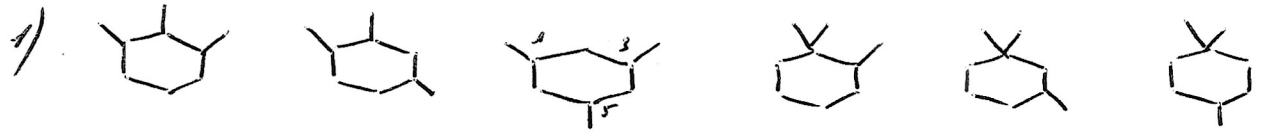
c) A cause de la liaison hydrogène intramoléculaire



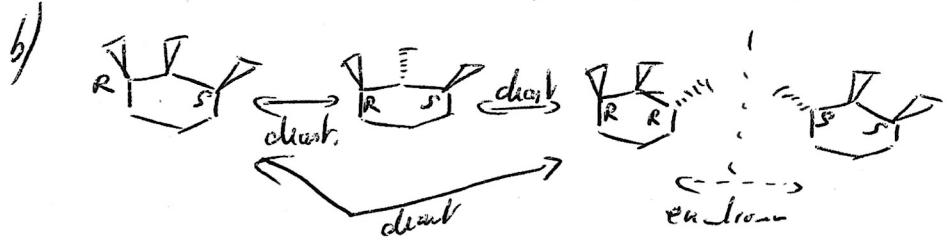
ou en Newman



Ex 11



Il n'y a pas de C* mais c'est de la pseudoasymétrie



en l'occurrence