



# Aromatic Compounds - II

Course on General Organic Chemistry for Class XI



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# IIT - ORGANIC CHEMISTRY

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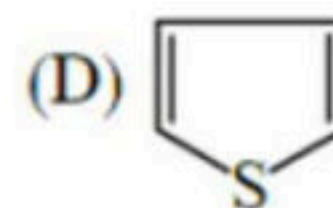
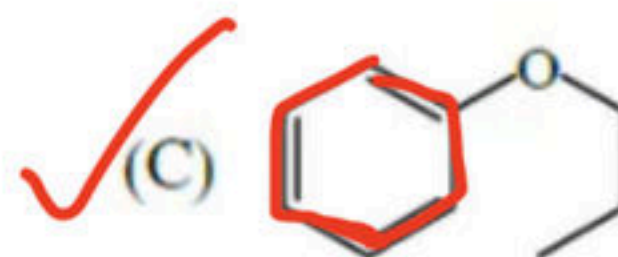
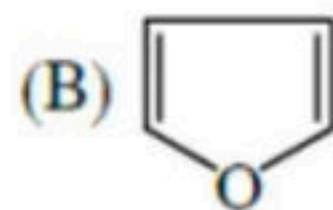
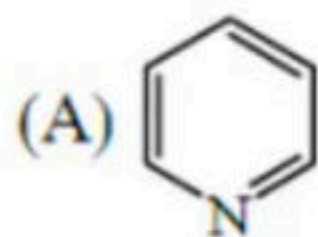
## NURTURE

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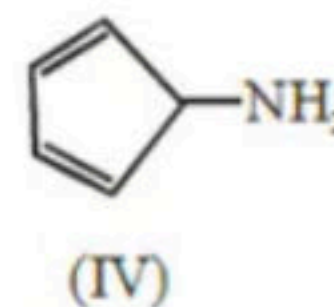
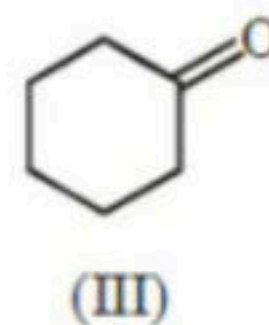
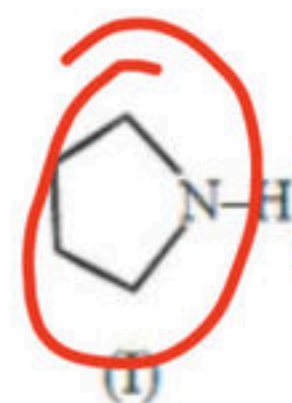
**Corporate Office: NAIVEDHYAM, Plot No. SP-11, Old INOX, Indra Vihar,  
Kota (Raj.) 324005**



1. **Statement 1 :** Urea is an organic compound. [3]  
**Statement 2 :** It can be synthesized only by living organism.
- (A) Statement-1 is true, statement-2 is true and statement-2 is correct explanation for statement-1.  
 (B) Statement-1 is true, statement-2 is true and statement-2 is NOT the correct explanation for statement-1.  
 ✓ (C) Statement-1 is true, statement-2 is false.  
 (D) Statement-1 is false, statement-2 is true.
2. Compound which is not heterocyclic - [3]



Paragraph for Q. 3 to 4





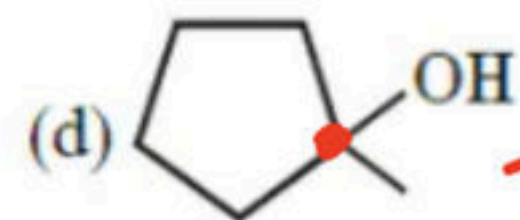
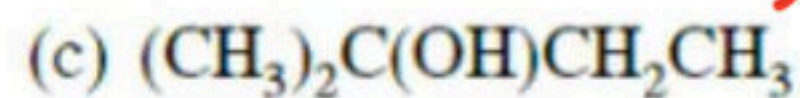
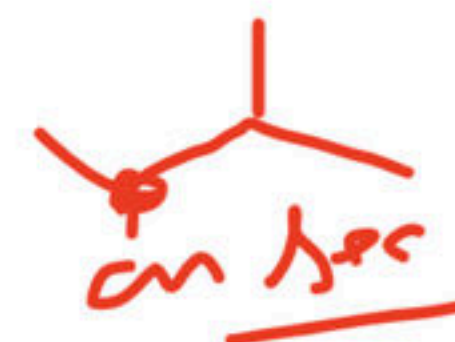
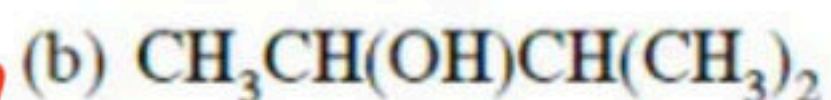
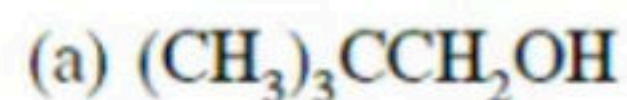
3. Heterocyclic compound is – [3]

- ✓ (A) I & II (B) II & III (C) III & IV (D) I, II, III & IV

4. Compound which is not amine – [3]

- (A) I (B) II ✓ (C) III (D) IV

5. Classify the following alcohols as primary, secondary, or tertiary: [3]

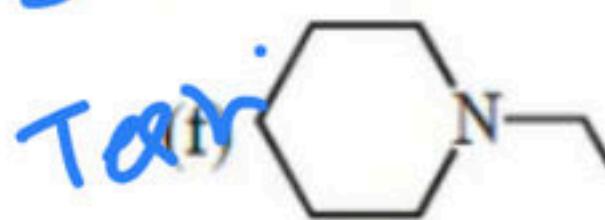
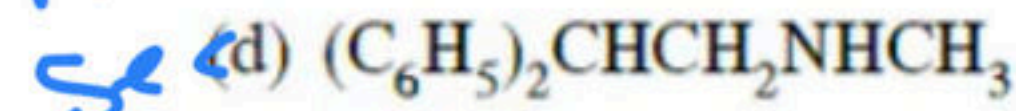
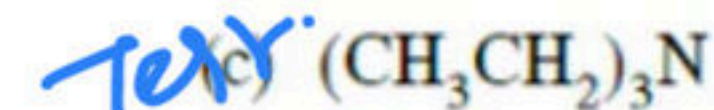
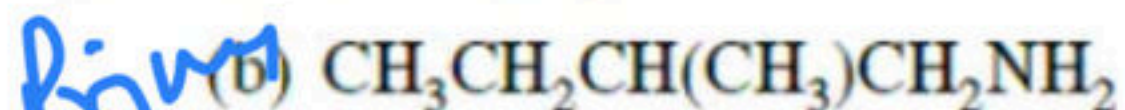


tertiary.



secondary.

6. Classify the following amines as primary, secondary, or tertiary: [3]






7. Which of the following have both 2° alcohol & 2° carbon only . [3]

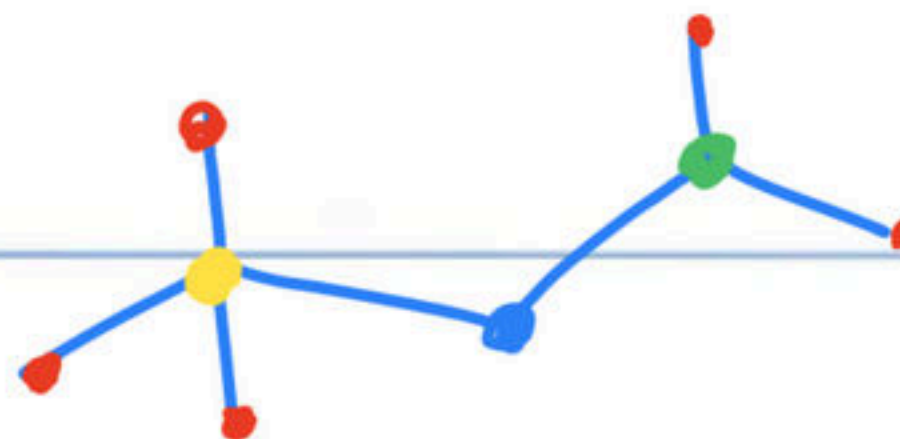


8. **Statement 1 :** If number of  $\pi$  bonds in the compound is 3 then its degree of unsaturation must be 3  
*Because* [3]

**Statement 2 :** For one  $\pi$  bond degree of unsaturation is equal to 1

- (A) Statement-1 is true, statement-2 is true and statement-2 is correct explanation for statement-1.  
(B) Statement-1 is true, statement-2 is true and statement-2 is NOT the correct explanation for statement-1.  
(C) Statement-1 is true, statement-2 is false.  
 (D) Statement-1 is false, statement-2 is true.





9. Isooctane contains

[3]

(A) five ( $1^\circ$  -C), one ( $2^\circ$  -C), two ( $3^\circ$  -C) atoms

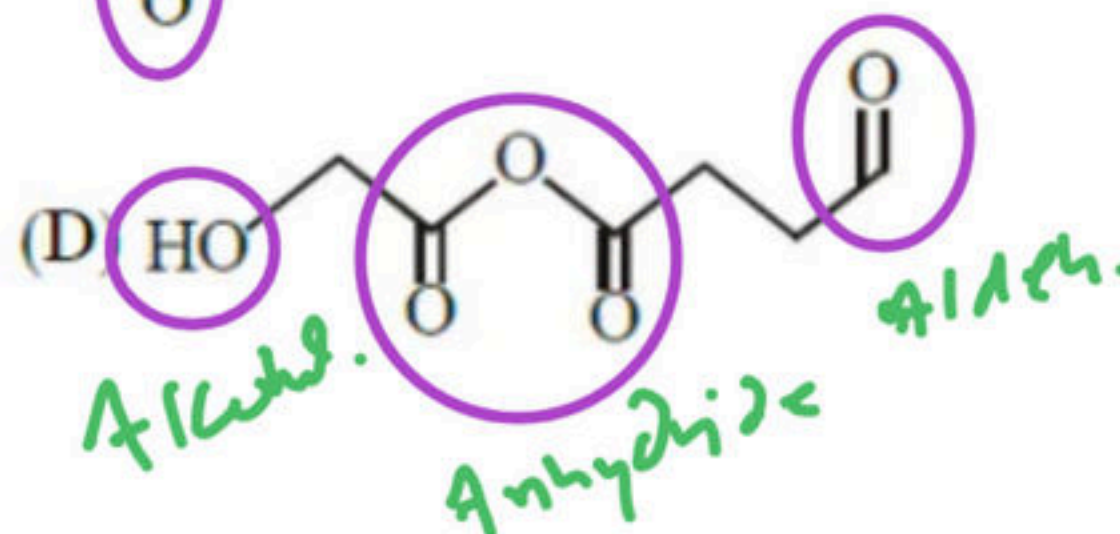
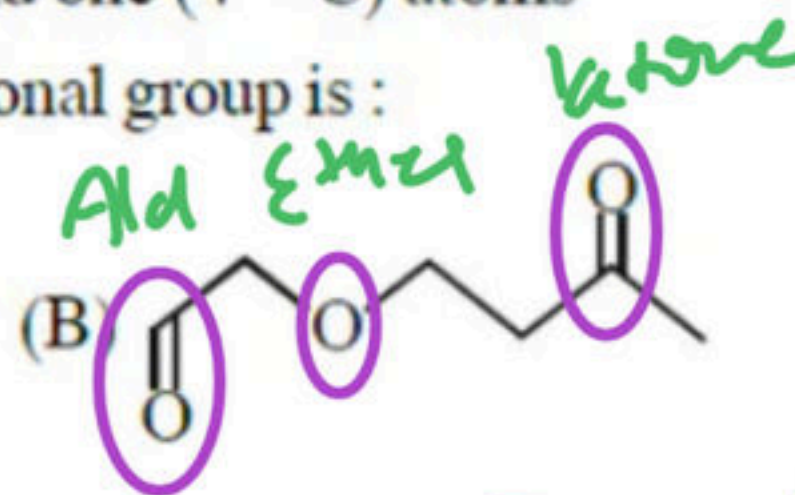
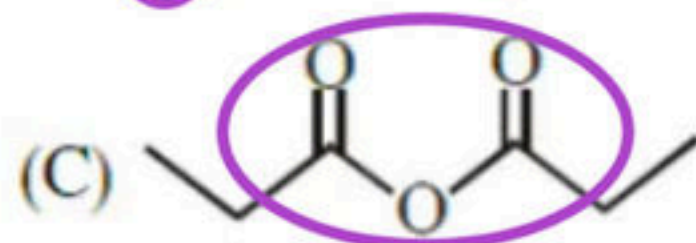
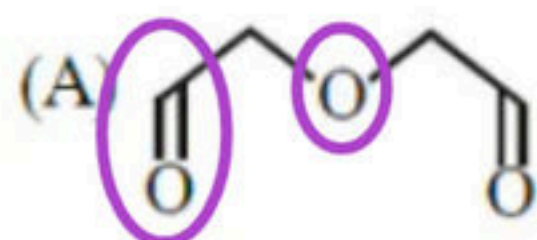
~~(B) four ( $1^\circ$  -C), two ( $2^\circ$  -C), one ( $3^\circ$  -C) and one ( $4^\circ$  -C) atoms~~

~~(C) four ( $1^\circ$  -C), two ( $2^\circ$  -C) and one ( $3^\circ$  -C) atoms~~

☒ (D) five ( $1^\circ$  -C), one ( $2^\circ$  -C), one ( $3^\circ$  -C) and one ( $4^\circ$  -C) atoms

10. Compound having only three different functional group is :

[4]



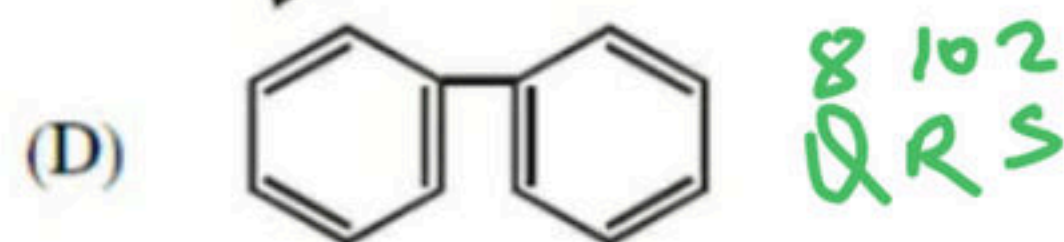
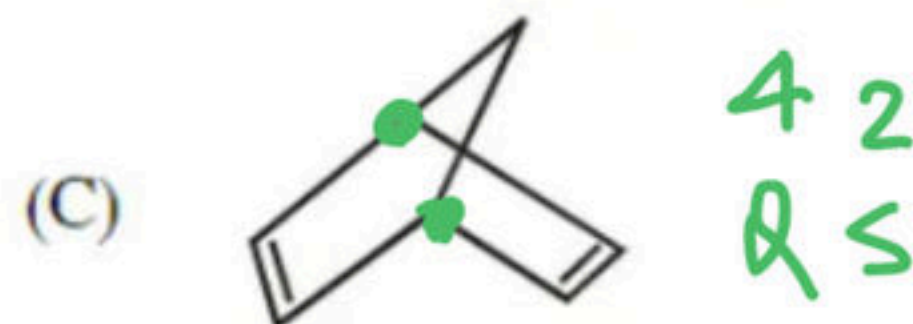
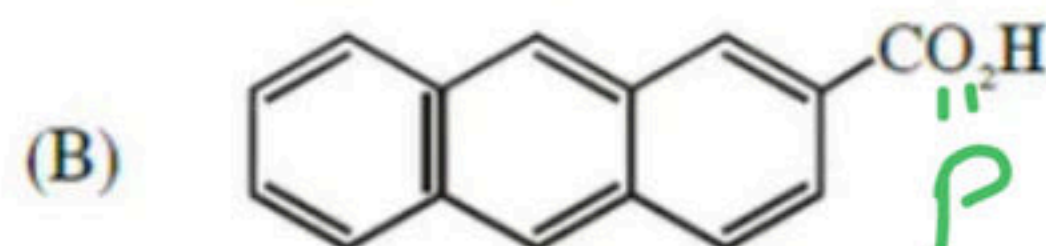
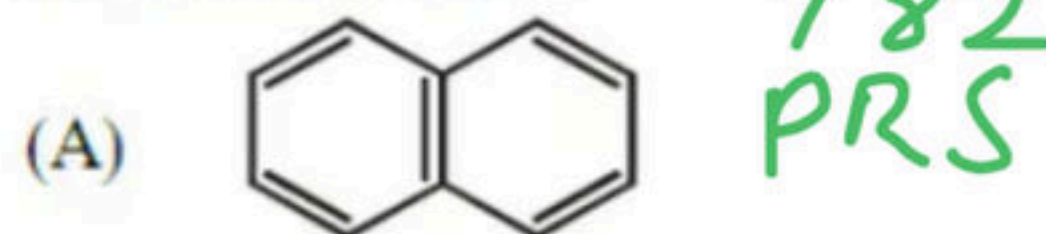


11. Match the column

[4]

Column I

(General formula)



Column II

(P) Index of hydrogen deficiency is odd

(Q) Index of hydrogen deficiency is Even

(R) Even number of 2° Carbon

(S) Even number of 3° Carbon







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## NURTURE

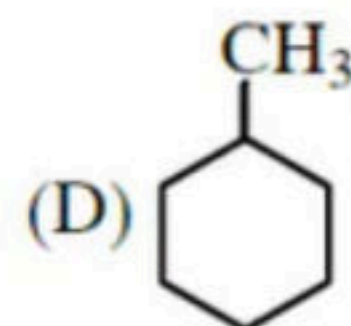
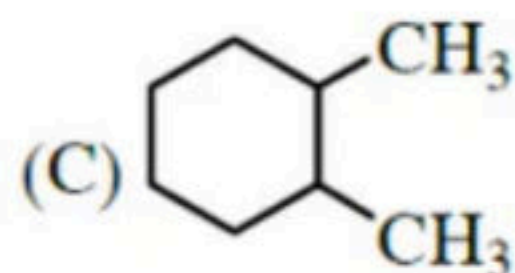
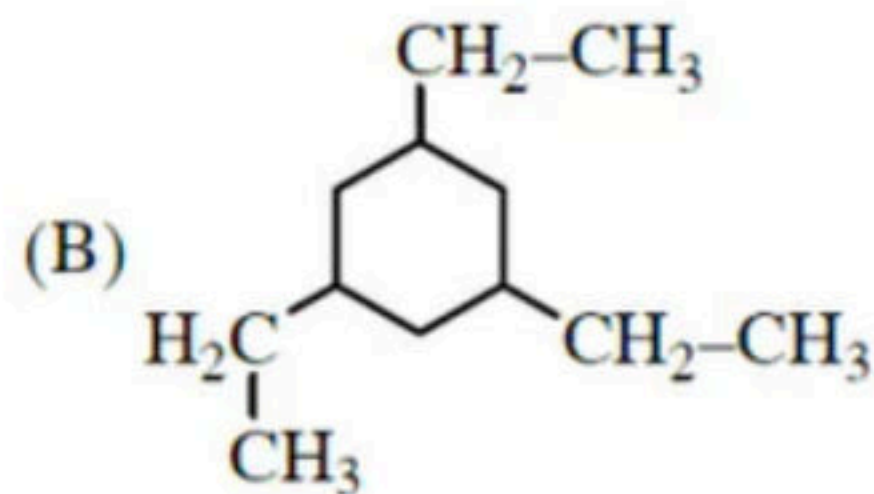
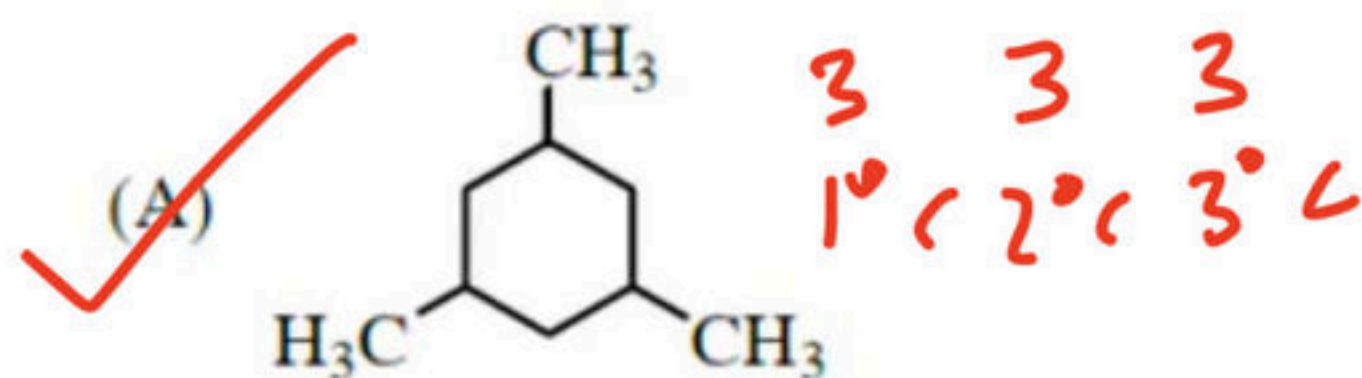
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1. **Statement 1 :** Phenol is a heterocyclic compound. [3]  
**Statement 2 :** In heterocyclic compound different atoms like O, N, S etc. are present in the ring.  
 (A) Statement-1 is true, statement-2 is true and statement-2 is correct explanation for statement-1.  
 (B) Statement-1 is true, statement-2 is true and statement-2 is NOT the correct explanation for statement-1.  
 (C) Statement-1 is true, statement-2 is false.  
 ✓ (D) Statement-1 is false, statement-2 is true.
2. In which compound  $1^\circ\text{C} : 2^\circ\text{C} : 3^\circ\text{C}$  (carbon) = 1 : 1 : 1.? [3]





{ Empir }

=

alkene

3. All the members of homologous series have same :

(A) molecular mass

(B) functional group

(C) empirical formula

(D) general molecular formula

4. Compound having molecular formula  $C_nH_{2n-4}O_3$  can have functional group.

(A) 3-Aldehyde group

(B) 1-Carboxylic acid & 2-Aldehyde

(C) 1-Carboxylic acid anhydride & 1-alcohol

(D) 1-Carboxylic acid & 1-alcohol

$DBE=3$

A & D

Emp. F  $CH_2$

$MF = (CH_2)^x$   
Alkene

$x=2$   $C_2H_4$  mf

$x=3$   $C_3H_6$  mf

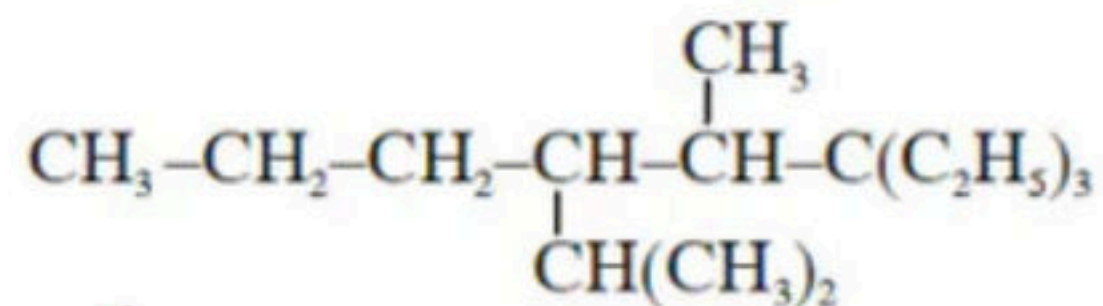
$x=4$   $C_4H_8$  mf

$n=2$



5. The correct IUPAC name of the compound is :

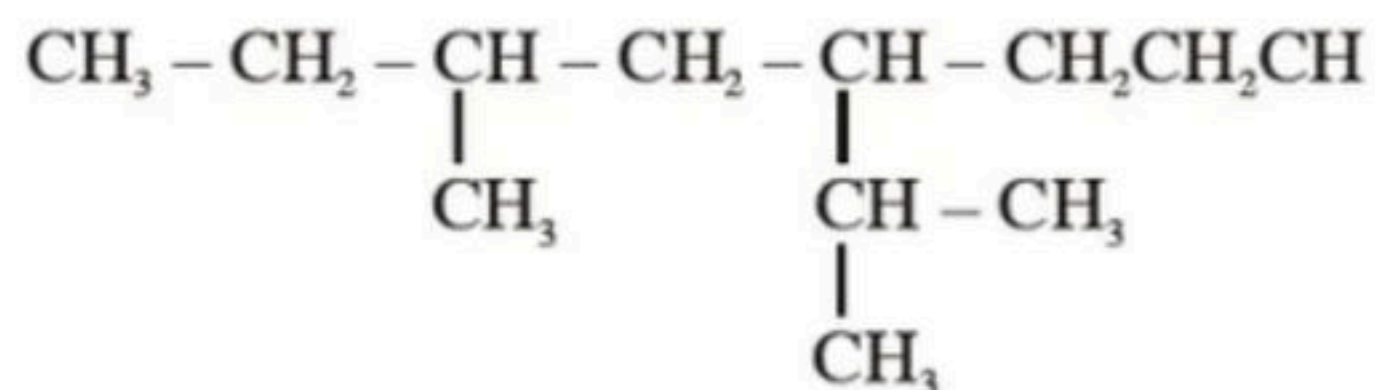
[3]



- ✓ (A) 3,3-Diethyl-4-methyl-5-(1-methylethyl) octane  
(B) 6,6-Diethyl-4-methyl-5-isopropyloctane  
(C) 6,6-Diethyl-3-methyl-5-(1-methylethyl) octane  
(D) 6,6-Diethyl-4-isopropyl-5-methyloctane

6. IUPAC name of the compound

[3]



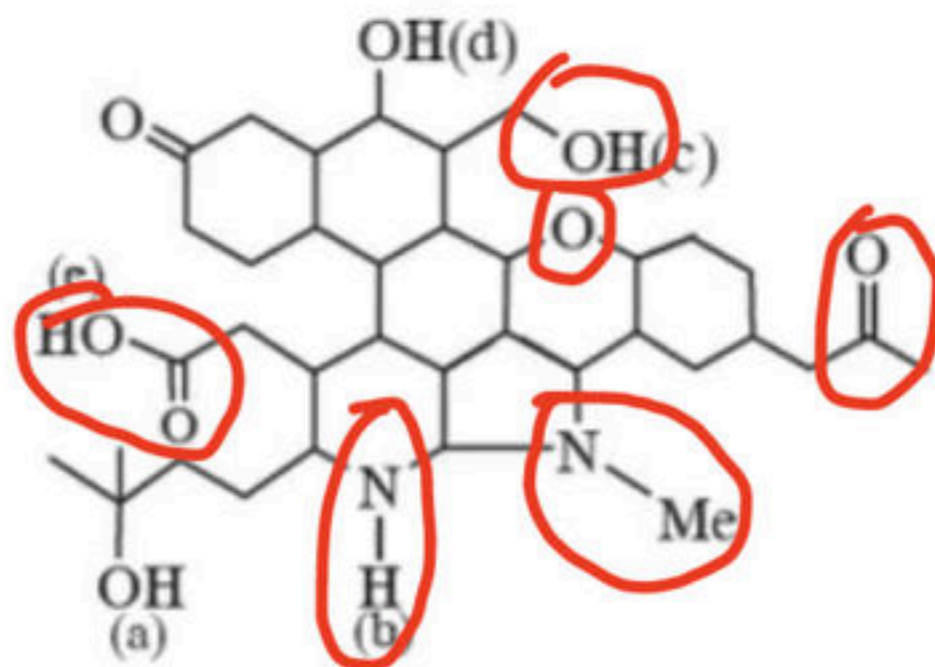
- (A) 4-Isopropyl-6-methyloctane  
(C) 3-Methyl-5-isopropyloctane  
✓ (B) 3-Methyl-5-(1-methylethyl) octane  
(D) 6-Methyl-4-(1-methylethyl) octane



7. IUPAC name of pivalic acid  $\left( \begin{array}{c} \text{H}_3\text{C} \\ \text{H}_3\text{C} \\ \text{H}_3\text{C} \end{array} \right) \text{C}-\text{COOH}$  is : [3]
- (A) Isobutylic acid (B) 2-carboxy-2-methyl propane  
 ✓ (C) 2,2-dimethyl propanoic acid (D) 2,2,2 trimethylethanoic acid

### Paragraph for Question 08 and 09

Observe following compound and answer questions given below :



8. Total number of different types of functional groups in this compound are [3]

(A) 5

✓ (B) 6

(C) 7

(D) 8

9. Degree of unsaturation of this compound is [3]

(A) 8

(B) 9

✓ (C) 10

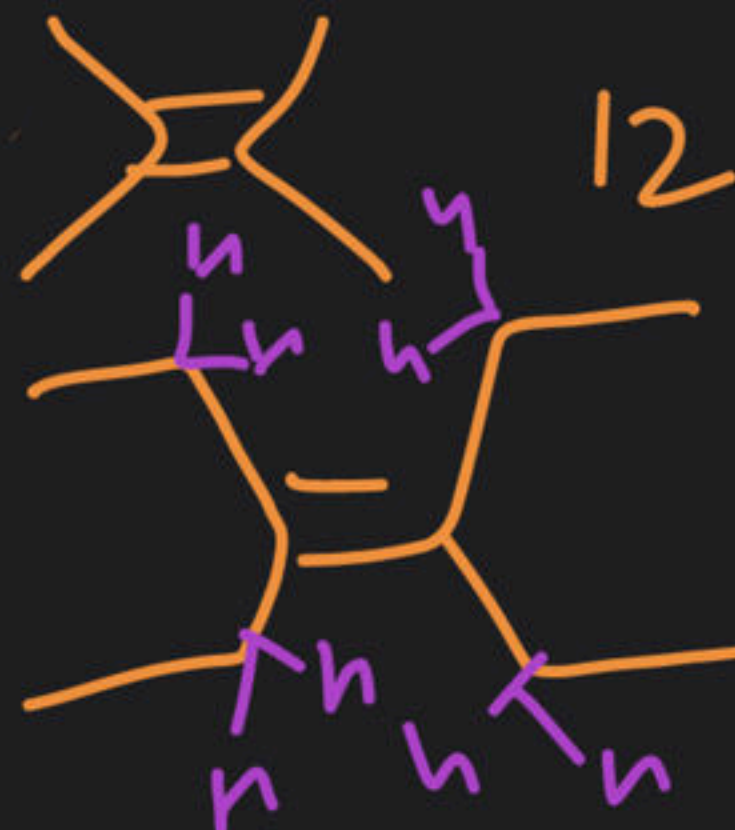
(D) 11

DPP  $\Rightarrow$  445



Ex-3:

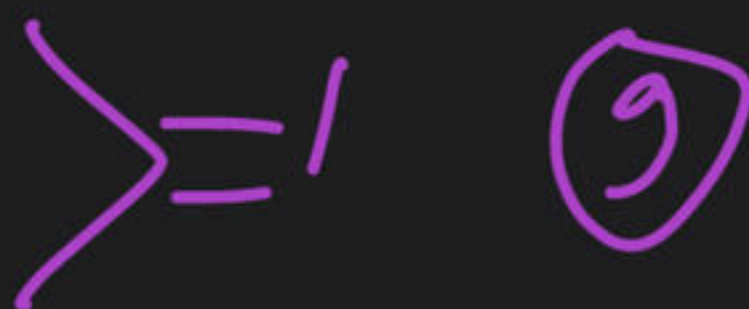
(vi)



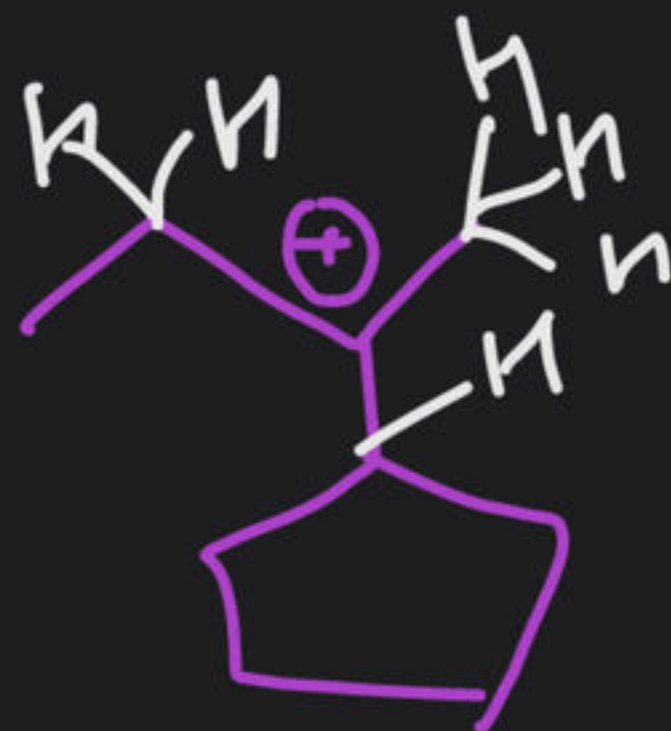
(8)

(vi)

(vii)



(viii)



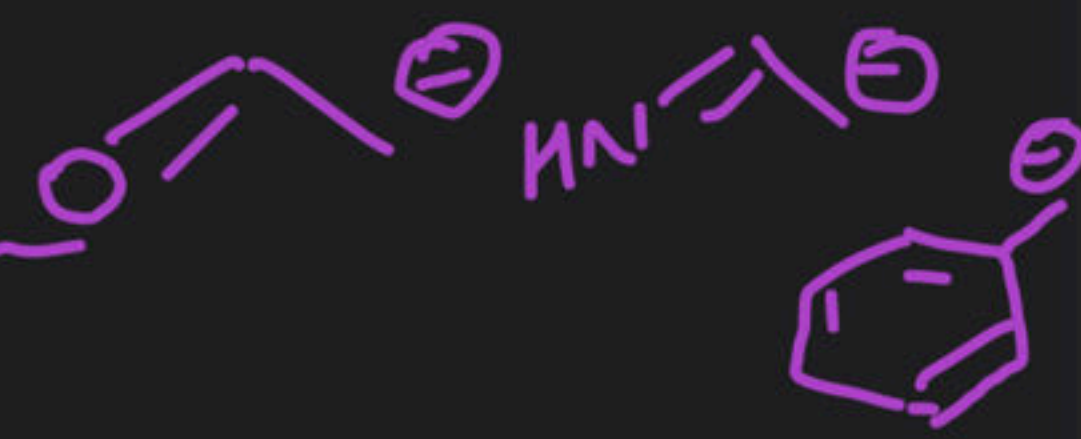
KE!

(3)

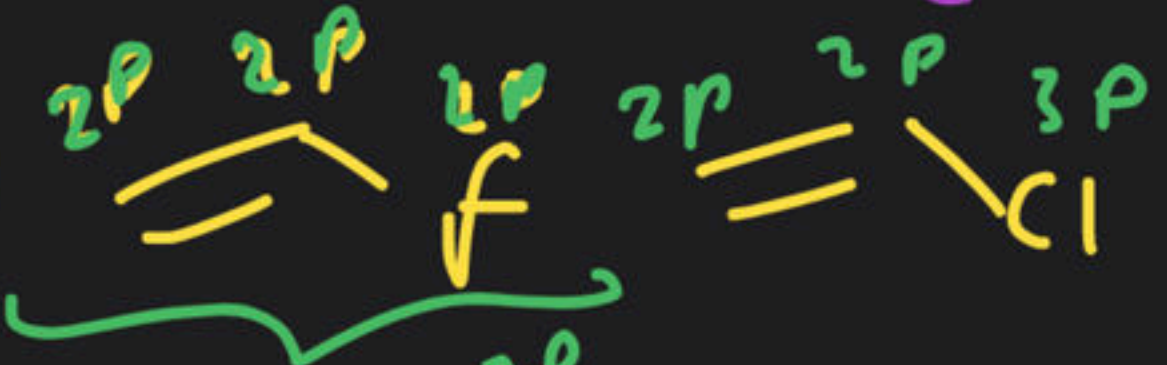
2 > 1

(4)

1 > 2 > 3

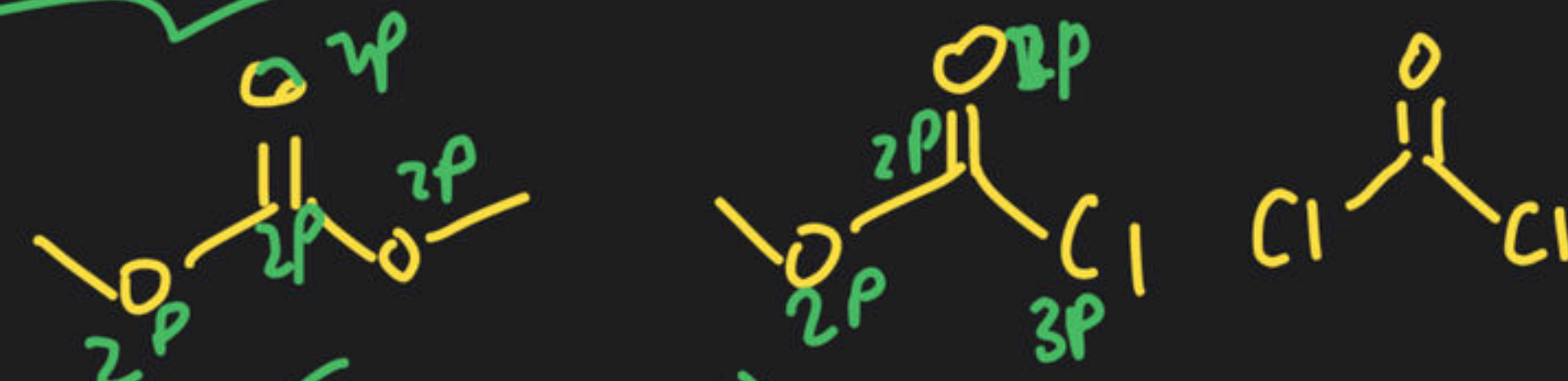


(5)



172

(6)



(1 > 2 > 3)

(7)

2 > 1

6p

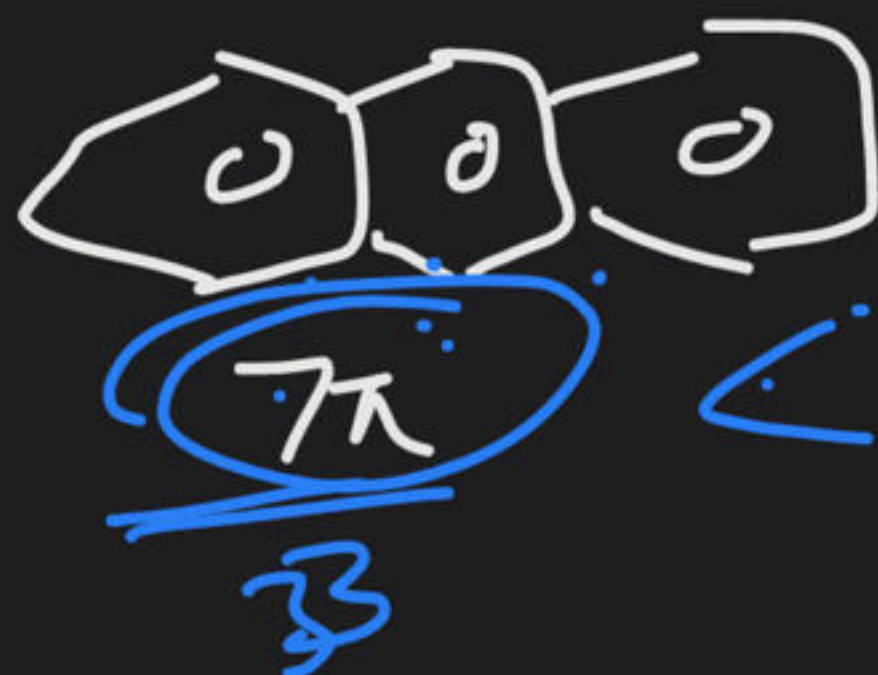
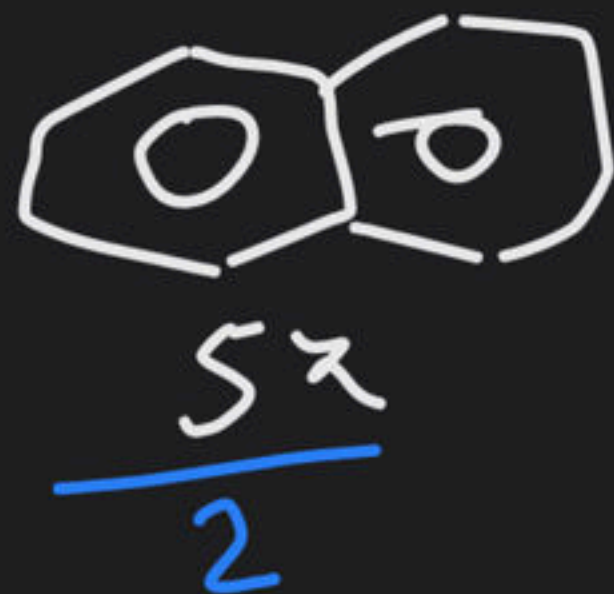
7p

(8)

4 > 3 > 2 > 1



(9)



$1 > 2 > 4 > 3$

(10)  $1 > 2$  (SZR effect)

(11)  $1 > 2$  (extended > cons)

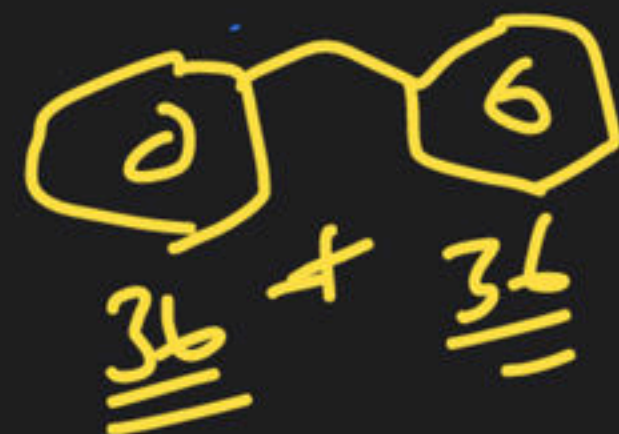
(12)  $2 > 1$  (ε > cons)

(13)  $1 > 2$  (ε > cons)

(14)  $(1-2)$

(15)

(16)  $2 > 1$

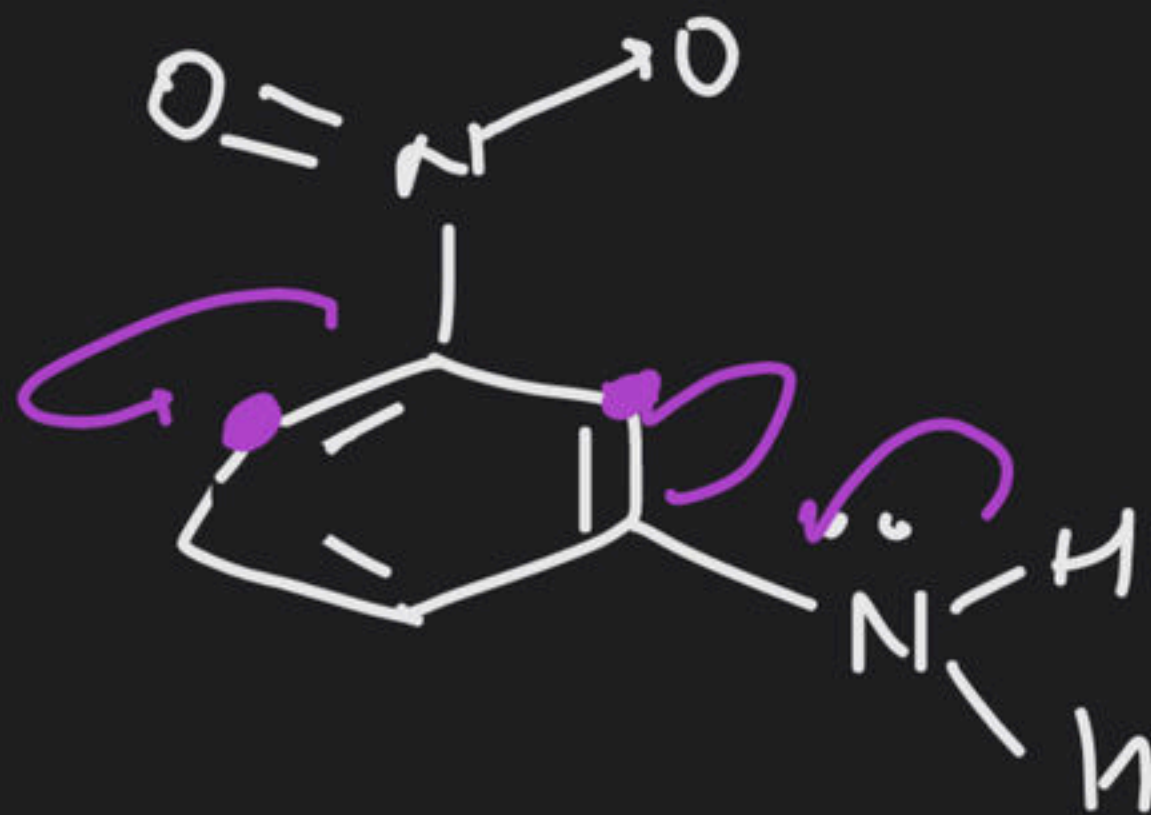
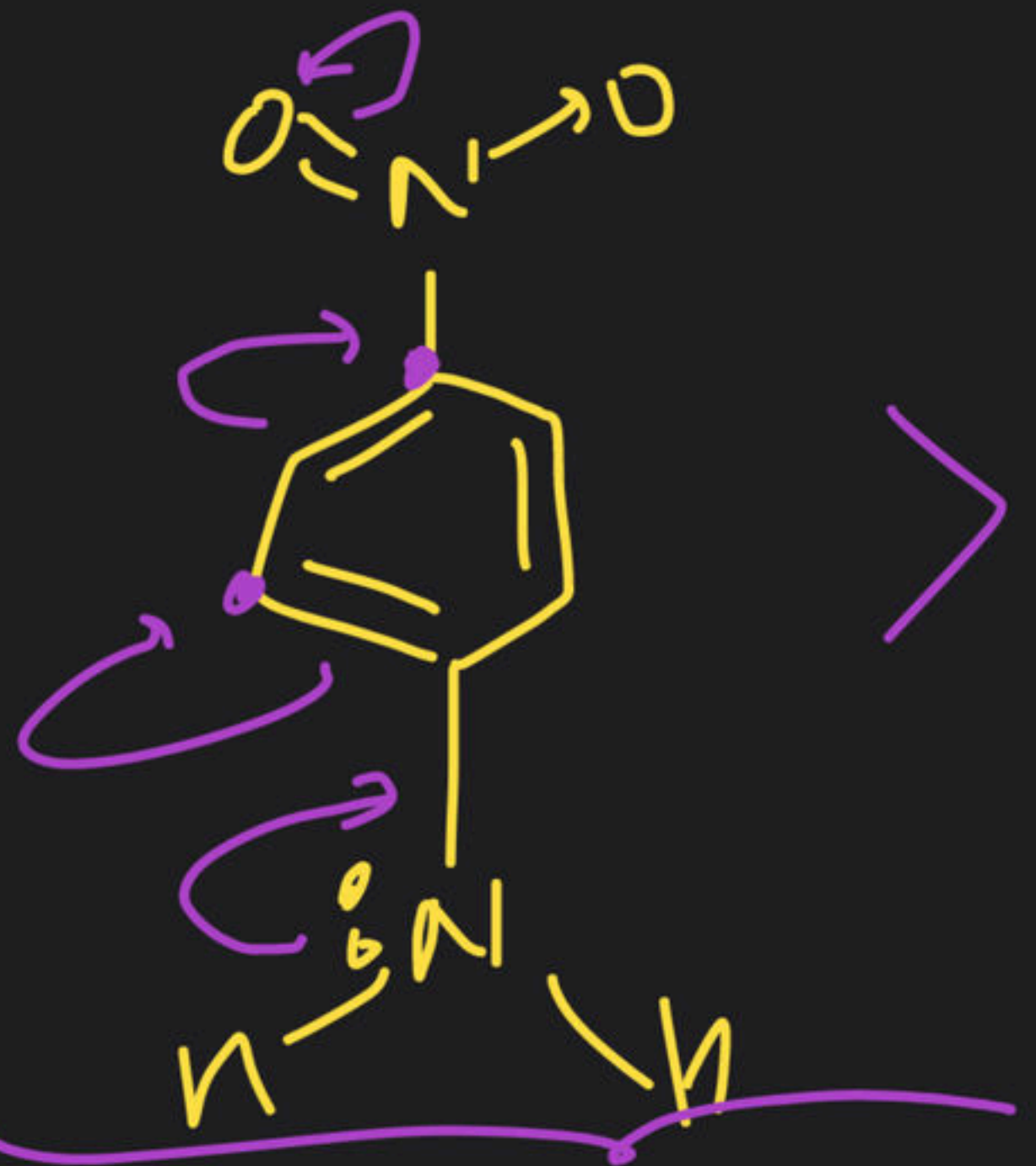


(17)

$4 > 3 > 2 > 1$

(21)  $1 > 2$





(6)



17273

NON

(3)

271

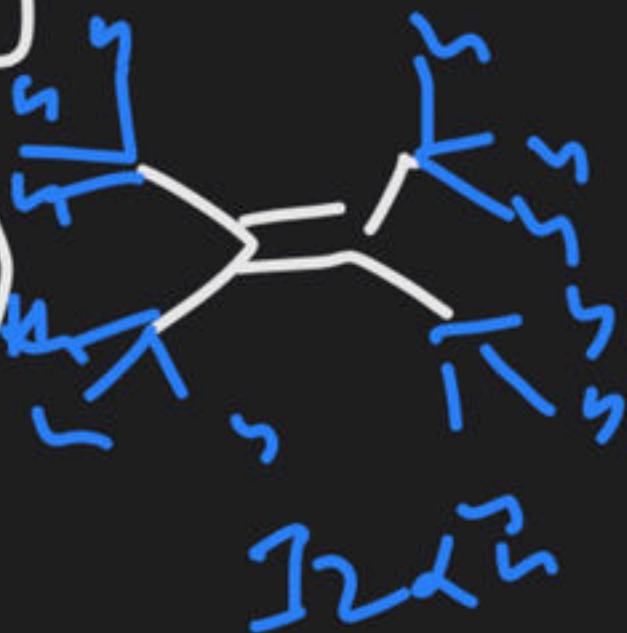
(4)

17372  $\Rightarrow$  (1/stability)

(5)

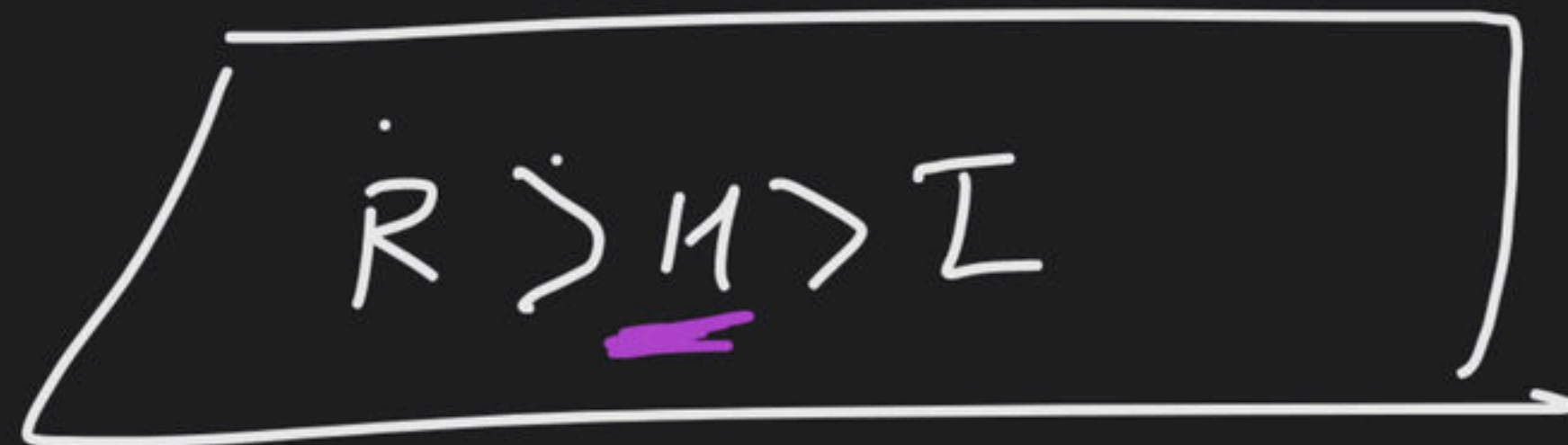
27371

55 51 29



$\frac{51}{3}$

(7)  $3 > 2 > 1$



(8)  $1 > 2$   
 Resonance  
 Stability  $\left( \frac{1}{\text{Stability}} \right)$

~~Imp~~  
 (9)



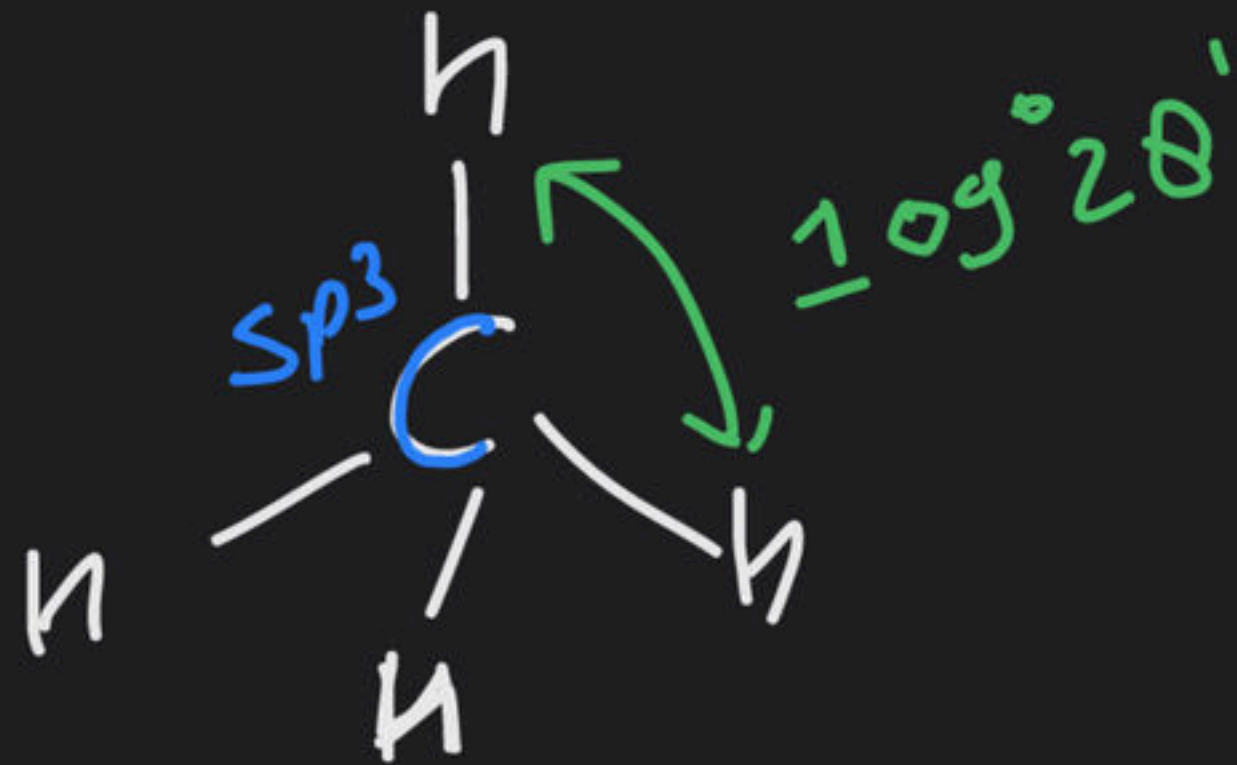
$\left( \frac{1}{\text{Stability}} \right)$

(10)  $2 > 1$   $\left( \frac{1}{\text{Stability}} \right)$



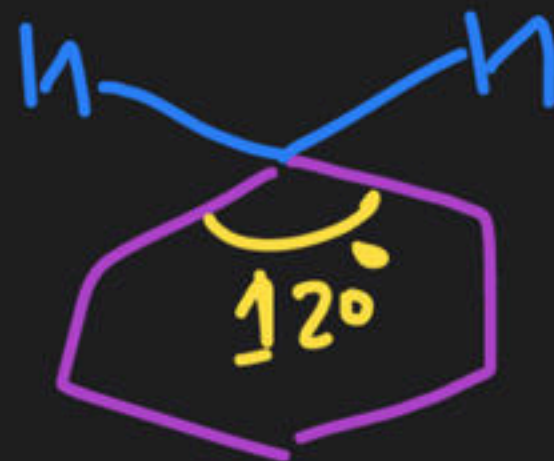


## (#) Bayer's Strain Angle theory



For any  $sp^3$  hybridised Carbon atom <sup>Bond</sup> angle must be  $109^{\circ}28'$  for maximum Stability

For following Cycloalkanes (it was believed till that cycloalkanes are planar).

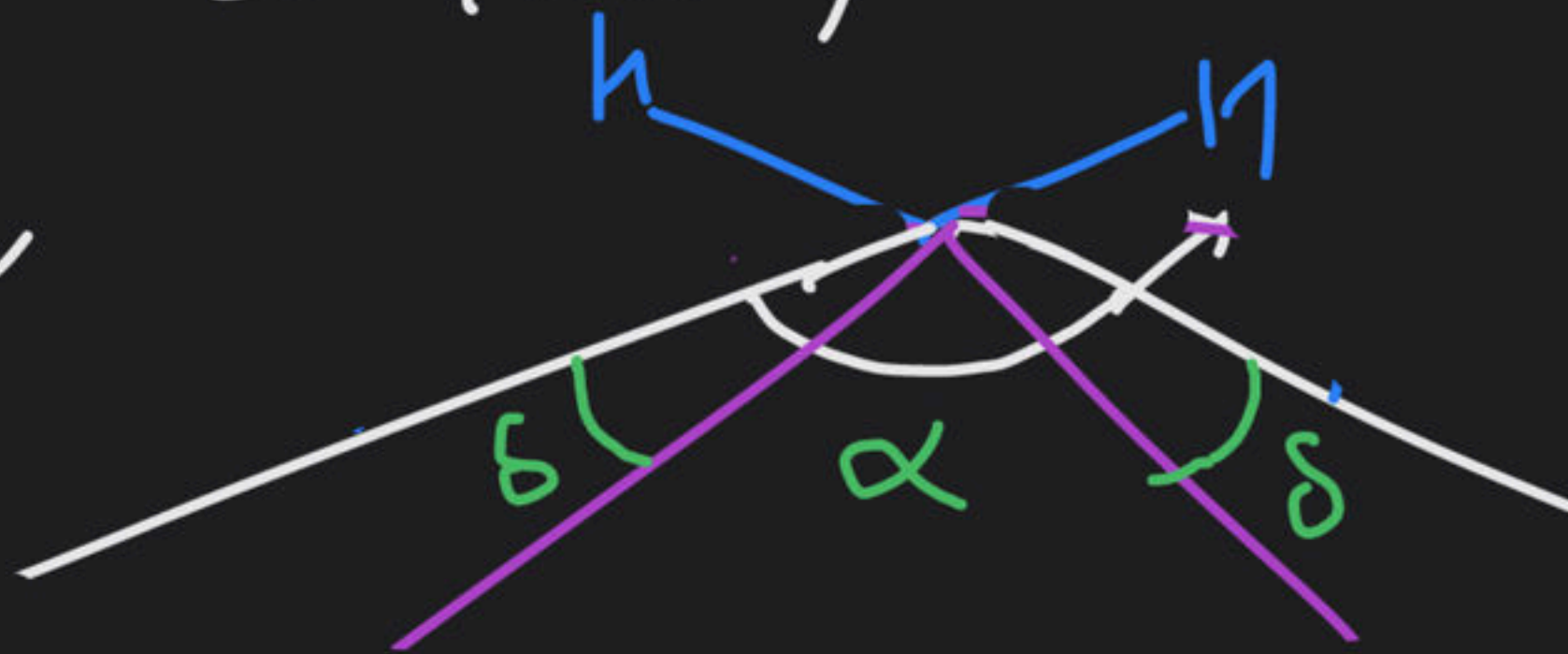


Acc. to Baeyer:

$\Rightarrow$  All cyclo alkanes are strained & are not stable

$\Rightarrow$  Strain  $\propto$   $\frac{1}{\text{Stability}}$

$\Rightarrow$  Acc. to Baeyer





$$\Rightarrow S + \alpha + \delta = 109^\circ 28'$$

$$\Rightarrow 2\delta + \alpha = 109^\circ 28'$$

$$\Rightarrow 2\delta = 109^\circ 28' - \alpha$$

$$\Rightarrow \delta = \frac{1}{2} [109^\circ 28' - \alpha]$$

Strain in Cycloalkenes:

$$\underline{\alpha = 6^\circ} \Rightarrow \delta = +24.5^\circ$$

$$\underline{\alpha = 9^\circ} \Rightarrow \delta = +9.5^\circ$$

$$\underline{\alpha = 100^\circ} \Rightarrow \delta = 0.5^\circ$$

$$\underline{\alpha = 120^\circ} \Rightarrow \delta = -5.5^\circ$$

Acc. to Baeyer's

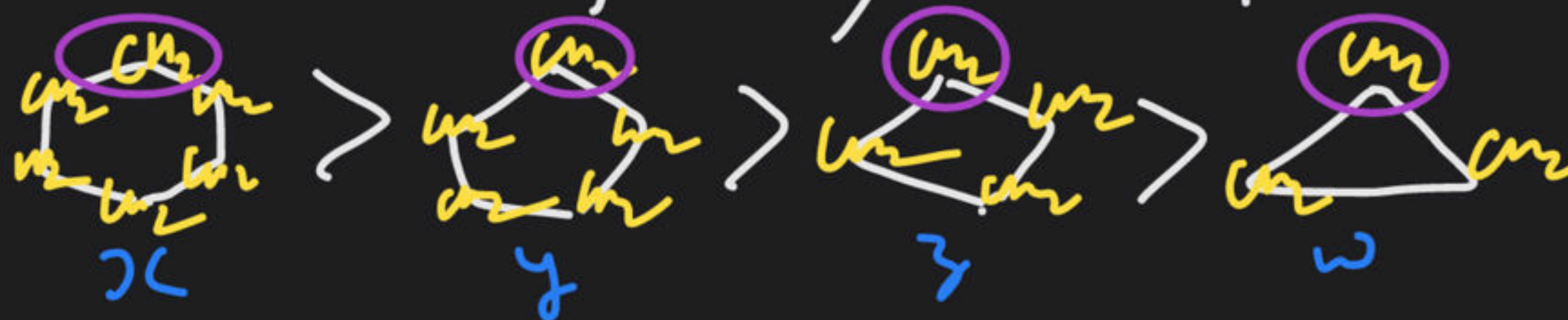
(\*) Strain order in cycloalkane



(\*) Stability order



⇒ But HOC data for cycloalkanes is





HOC per CH<sub>2</sub> for cycloalkanes (Strain)

Actual



Stability order

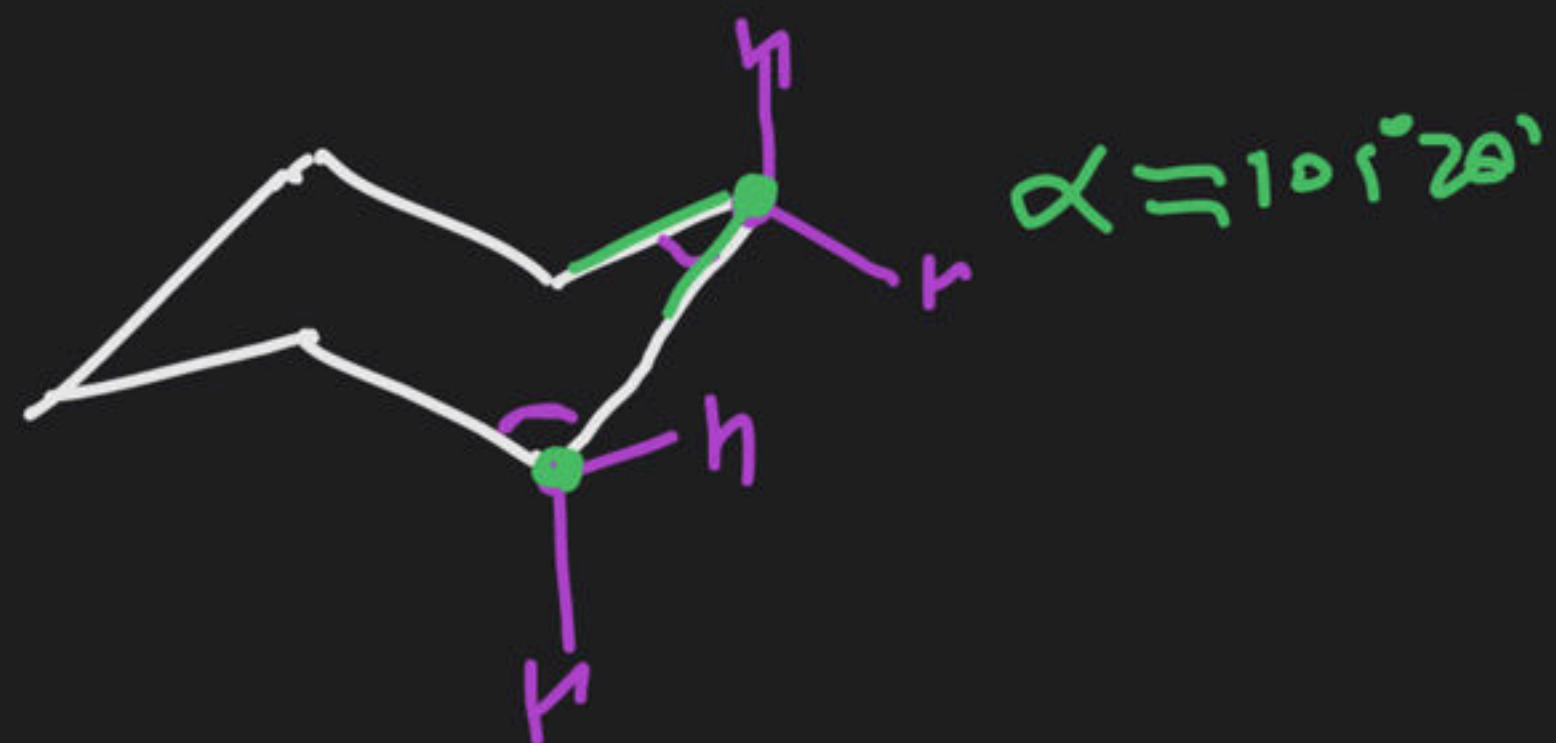
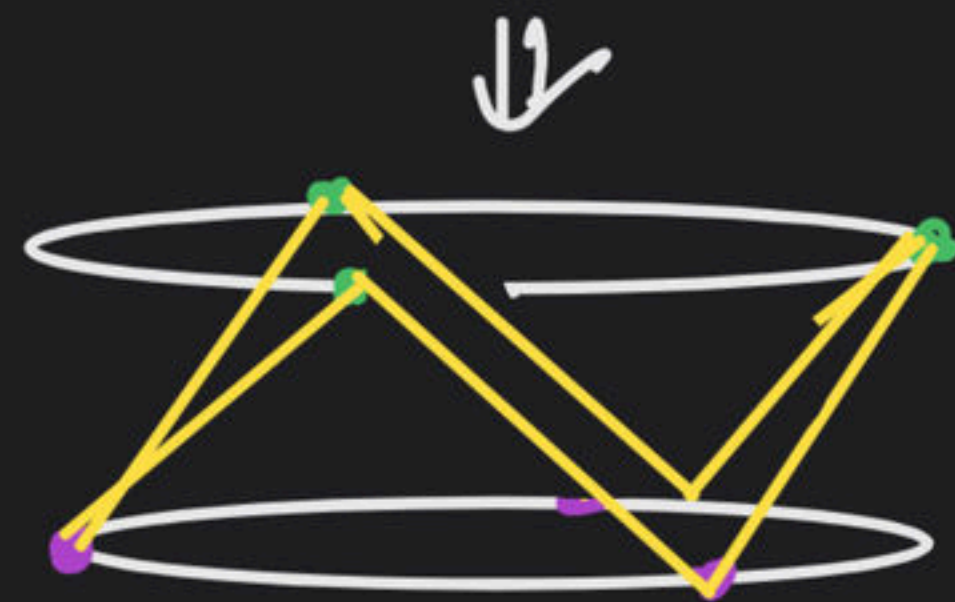


0.5°

-5.5°

It can be explain by understanding that cycloalkanes are not planar (except cyclopropane)

For cyclohexane



$(S \approx 0)$



(#) Heat of Combustion (HOC)

$\Rightarrow$  Enthalpy change when 1 mole of any compound is completely burnt/oxidised.

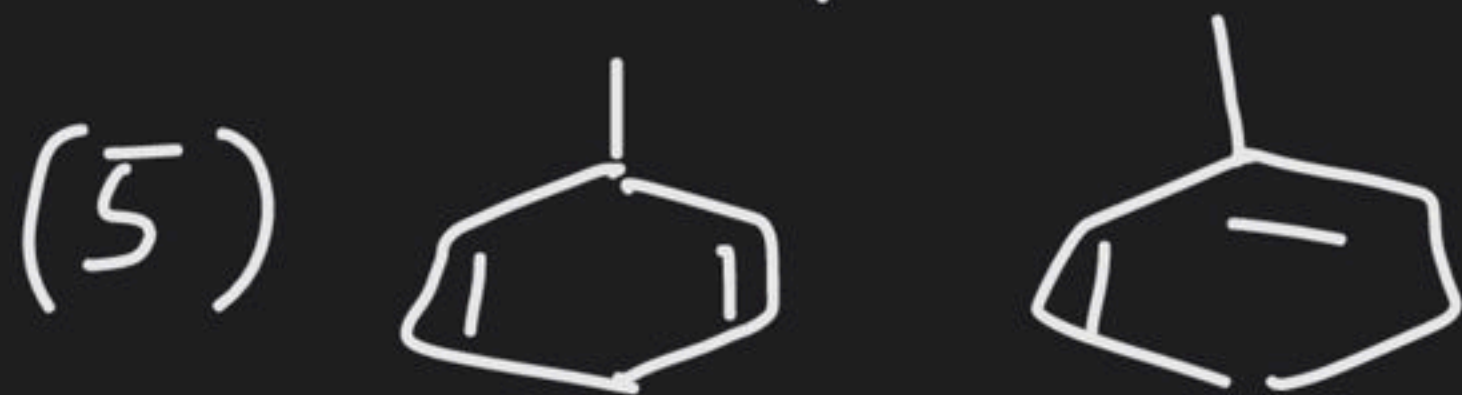
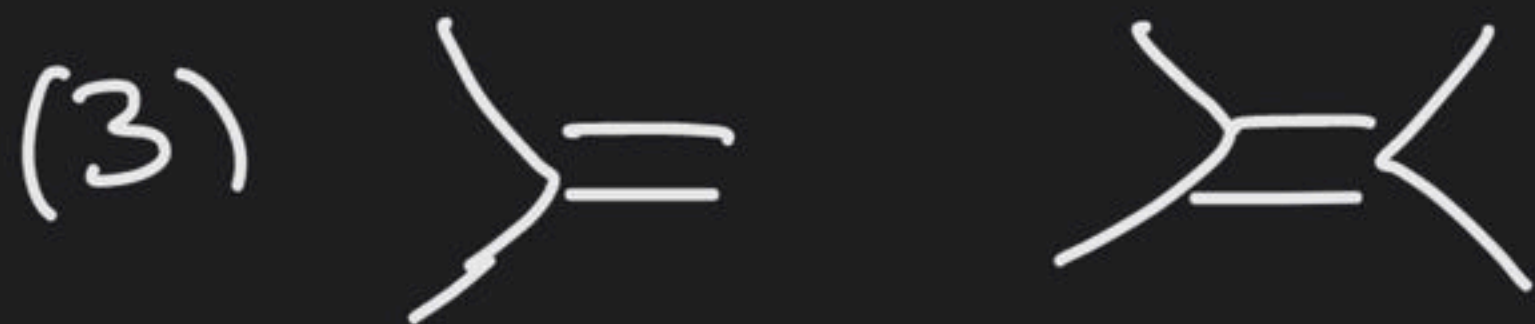
$$\text{HOC} \propto \text{No. of Carbon atom}$$

$\boxed{\text{HOC}/n_2} \propto \frac{1}{\text{Stability}} \propto \text{strain}$

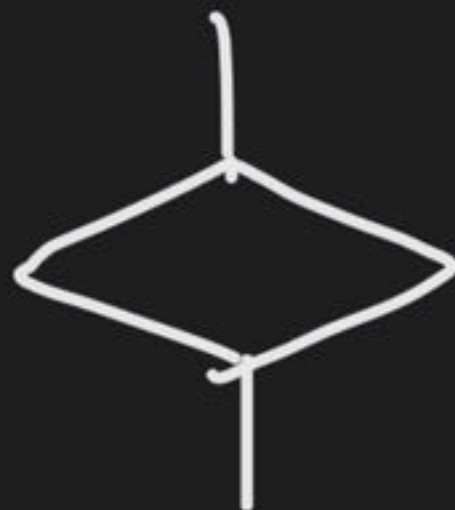
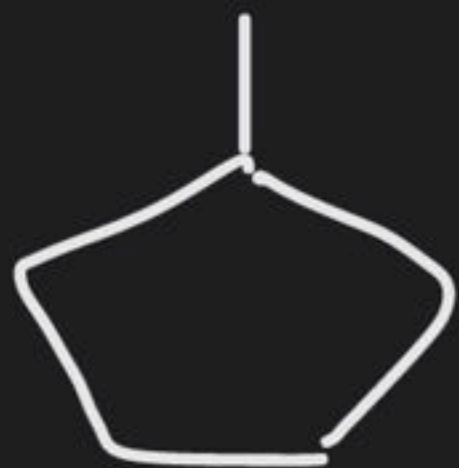
Note Actual stability order of compounds is decided by HOC.



(2) 2-methyl-3-ethyl pentane , 2,2,3 Trimethyl Butane





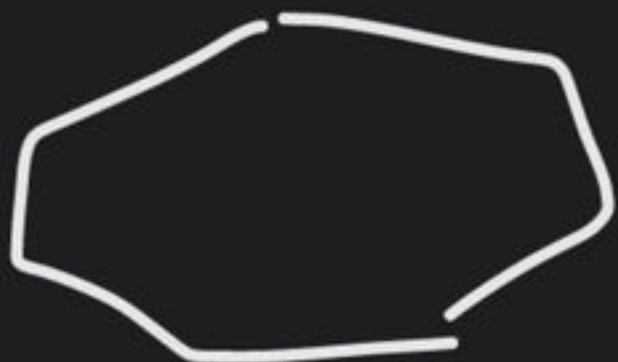


(7)

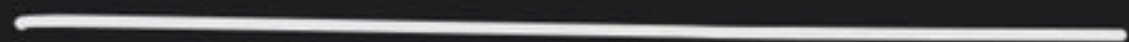


(8) - - - - - per  $\text{cm}_2$

(g)



(17)



for  $U_2$











# (1) Stability of intermediates

(a) Stability of Carbocation  $\propto$  ED-groups



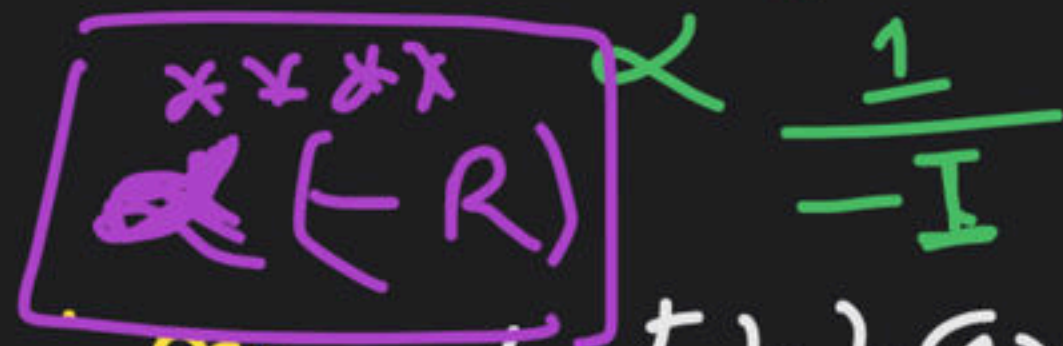
$$\propto (+R > +H > +I)$$

$$\propto \frac{1}{-I > -H > -R}$$

(b) Stability of Carbon free Radical  $\propto$  ED-group



$$\propto (+R > +H > +I)$$



(c) Stability of Carbanion  $\propto$  EW-groups



$$\propto (-R > -I)$$

$$\propto \frac{1}{(+I > +R)}$$





























