



GGSRDN

Educational Services Private Limited

9th, 10th, NEET, JEE (Main/Advanced)

अभ्यास ही सबसे बड़ा गुरु है।

CLASS : XI

(ORGANIC CHEMISTRY)

D P P

DAILY PRACTICE PROBLEM

DPP-1 to 16

ORGANIC CHEMISTRY
DPP
 DAILY PRACTICE PROBLEMS

DPP No. 1

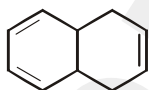
Total Marks : 32

Max. Time : 35 min.

Topic : IUPAC Nomenclature & Isomerism
Type of Questions

| Type of Questions | M.M., Min. |
|--|----------------------------|
| Single choice Objective ('-1' negative marking) Q.1 to Q.4 | (3 marks, 3 min.) [12, 12] |
| Multiple choice objective ('-1' negative marking) Q.5 to Q.6 | (4 marks, 4 min.) [8, 8] |
| Subjective Questions ('-1' negative marking) Q.7 | (4 marks, 5 min.) [4, 5] |
| Match the Following (no negative marking) Q.8 | (8 marks, 10 min.) [8, 10] |

1. Number of hydrogen atoms in the given compound is :



- (A) 8 (B) 10 (C) 12 (D) 14

 2. Ketene $\text{CH}_2 = \text{C} = \text{O}$ has

- (A) Only sp^2 carbon atom (B) Only sp carbon atom
 (C) sp^2 and sp carbon atoms (D) sp^3 , sp^2 and sp carbon atoms

3. Which of the following is not an unsaturated compound.

- (A) $\text{CH}_3 - \text{CH} = \text{CH} - \text{CH}_3$ (B) $\text{HC} \equiv \text{C} - \text{C} \equiv \text{CH}$
 (C) $\text{CH}_3 - \text{CH}_2 - \text{CH}_2 - \text{HC} \begin{matrix} \text{CH}_2 \\ | \\ \text{CH}_2 \end{matrix}$ (D) $\text{CH}_2 = \text{CH} - \text{CH} = \text{CH}_2$

4. Which of the following is homocyclic compound.



5.* The alicyclic compound/s is / are



6.* The correct options for a homologous series

- (A) All members have same general formula
 (B) All members have same chemical properties
 (C) All members have same physical properties
 (D) All members have same functional groups

Topic : IUPAC Nomenclature & Isomerism

Type of Questions

Single choice Objective ('-1' negative marking) Q.1 to Q.5

(3 marks, 3 min.)

M.M., Min.

[15, 15]

Multiple choice objective ('-1' negative marking) Q.6

(4 marks, 4 min.)

[4, 4]

Subjective Questions ('-1' negative marking) Q.7

(4 marks, 5 min.)

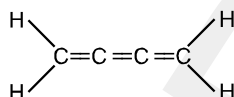
[4, 5]

Match the Following (no negative marking) Q. 8

(8 marks, 10 min.)

[8, 10]

1. Number of sp^2-sp sigma bonds in the given compound is :



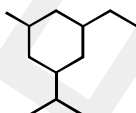
(A) 1

(B) 2

(C) 3

(D) 4

2. How many tertiary carbon atom are present in the compound :



(A) 2

(B) 3

(C) 4

(D) 5

3. Which statement is incorrect :

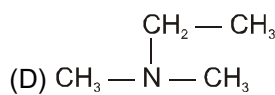
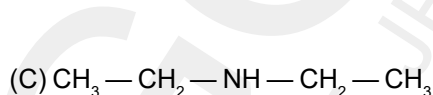
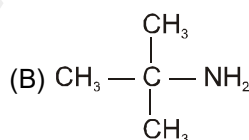
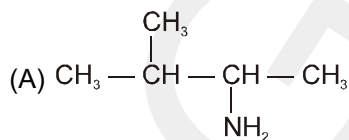
(A) $C_n H_{2n-2}$ is the general formula of alkyne

(B) $C_n H_{2n+2} O$ is the general formula of alkanol

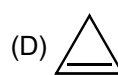
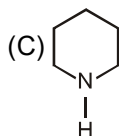
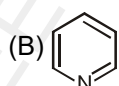
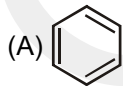
(C) $C_n H_{2n}$ is the general formula of alkene

(D) $C_n H_{2n+2}$ is the general formula of cycloalkane

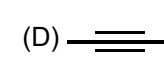
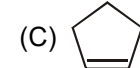
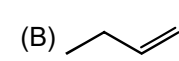
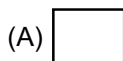
4. Which of the following is 3° -Amine :



5. Which of the following is heteroaromatic compound :



6.* Which of the following has $C_n H_{2n}$ general formula :



7. Calculate molecular weight of the lowest alkane containing a sequence of 1° , 2° , 3° and 4° carbon atoms.

8. Match the following :

| Column I | Column II |
|---------------------------|---------------------------|
| (A) 4 carbon atoms alkane | (P) Molecular weight = 26 |
| (B) 2 carbon atoms alkyne | (Q) Molecular weight = 42 |
| (C) 3 carbon atoms alkene | (R) Molecular weight = 40 |
| (D) 3 carbon atoms alkyne | (S) Molecular weight = 58 |

ORGANIC CHEMISTRY

DPP

DAILY PRACTICE PROBLEMS

DPP No. 3

Total Marks : 32

Max. Time : 35 min.

Topic : IUPAC Nomenclature & Isomerism

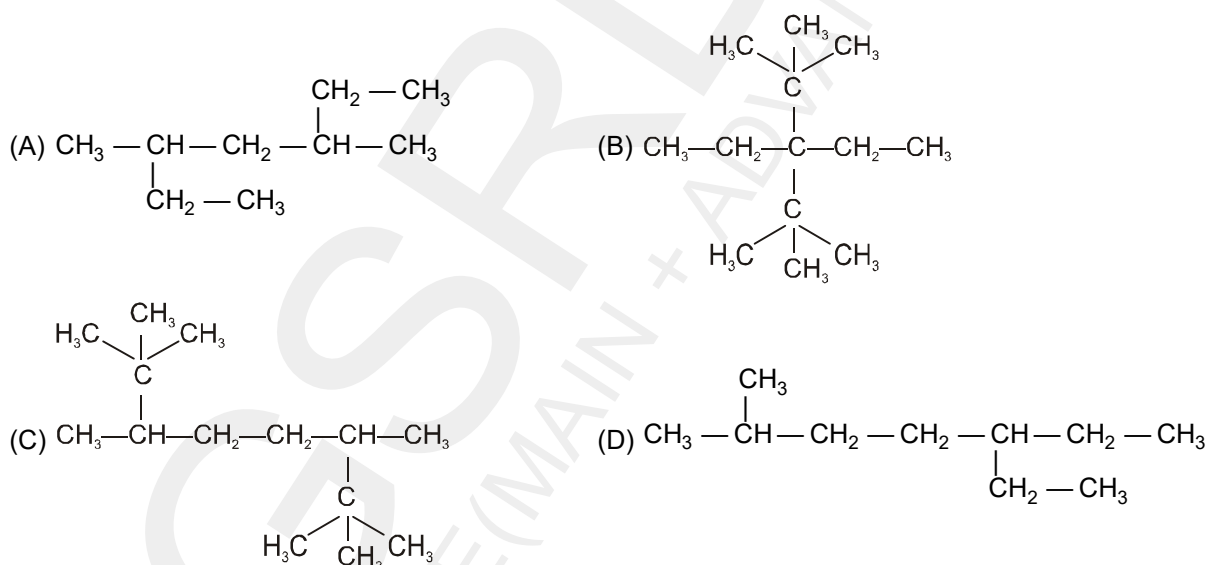
Type of Questions

| Type of Questions | M.M., Min. |
|--|----------------------------|
| Single choice Objective ('-1' negative marking) Q.1 to Q.4 | (3 marks, 3 min.) [12, 12] |
| Multiple choice objective ('-1' negative marking) Q.5 to Q.6 | (4 marks, 4 min.) [8, 8] |
| Subjective Questions ('-1' negative marking) Q.7 | (4 marks, 5 min.) [4, 5] |
| Match the Following (no negative marking) Q. 8 | (8 marks, 10 min.) [8, 10] |

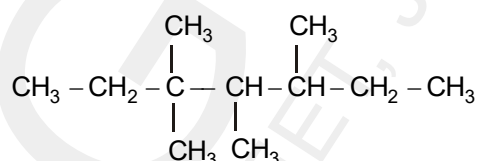
1. Which of the following alkanes contains primary, secondary, tertiary and quaternary carbon atoms together.

- (A) $(\text{CH}_3)_3\text{CH}$ (B) $(\text{C}_2\text{H}_5)_3\text{CH}$ (C) $(\text{CH}_3)_3\text{CCH}_2\text{CH}(\text{CH}_3)_2$ (D) $(\text{CH}_3)_4\text{C}$

2. Which of the following has longest chain of carbon :



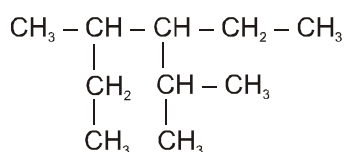
3. In following compound -



The correct lowest set of locants are

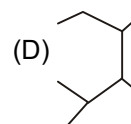
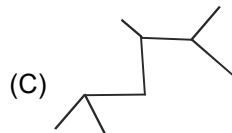
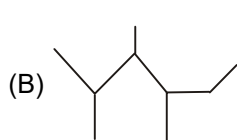
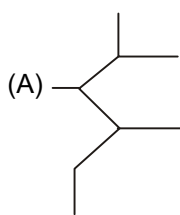
- (A) 3,3,4,5 (B) 3,4,5,5 (C) 4,5,3,3 (D) 5,5,4,3

4. The correct IUPAC name of the following compound is



- (A) 4-Ethyl-3,5-dimethylhexane (B) 2,4-Dimethyl-3-ethylhexane
 (C) 3-Isopropyl-4-methylhexane (D) 3-Ethyl-2,4-dimethylhexane

5.* The correct structure of 2,3,4-Trimethylhexane is :



6.* Choose the correct option's according to given IUPAC name :

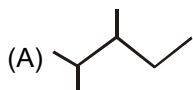
- (A) Neohexane is 2,2-Dimethylbutane.
 (B) Isobutane is 2-Methylpropane.
 (C) Isopentane is 2-Methylbutane.
 (D) Neopentane is Dimethylpropane.

7. Calculate the molecular weight of the lowest hydrocarbon which contains sp & sp^2 hybridised carbon atoms only.

8. Match the following :

Column-I

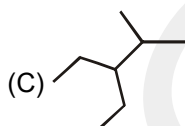
Column-II



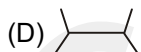
(p) 3-Ethyl-2-methylpentane



(q) 3-Methylhexane



(r) 2,3-Dimethylbutane



(s) 2,3-Dimethylpentane

ORGANIC CHEMISTRY
DPP

DAILY PRACTICE PROBLEMS

DPP No. 4

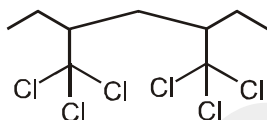
Total Marks : 27

Max. Time : 28 min.

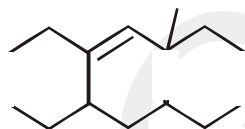
Topic : IUPAC Nomenclature & Isomerism
Type of Questions

| Type of Questions | M.M., Min. |
|--|----------------------------|
| Single choice Objective ('-1' negative marking) Q.1 to Q.4 | (3 marks, 3 min.) [12, 12] |
| Multiple choice objective ('-1' negative marking) Q.5 to Q.6 | (4 marks, 4 min.) [8, 8] |
| Subjective Questions ('-1' negative marking) Q.7 | (4 marks, 5 min.) [4, 5] |
| Comprehension ('-1' negative marking) Q.8 | (3 marks, 3 min.) [3, 3] |

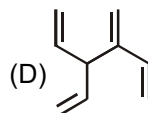
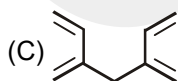
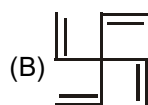
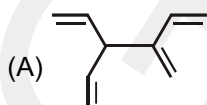
1. The correct IUPAC name of $\text{H}_3\text{C}-\overset{\text{Br}}{\underset{\text{F}}{\text{C}}}-\underset{\text{C}_2\text{H}_5}{\text{CH}}-\overset{\text{Br}}{\underset{\text{I}}{\text{C}}}-\text{CH}_3$
- (A) 2,4-Dibromo-3-ethyl-2-fluoro-4-iodopentane (B) 2,4-Dibromo-3-ethyl-4-fluoro-2-iodopentane
 (C) 2,4-Dibromo-4-fluoro-2-iodo-3-ethylpentane (D) 2,4-Dibromo-2-fluoro-4-iodo-3-ethylpentane
2. The correct IUPAC name of the following compound is :



- (A) 1,1,1,5,5,5-Hexachloro-2,4-diethylpentane. (B) 3,5-Hexachlorodimethylheptane.
 (C) 3,3,3,5,5,5-Hexachloromethylheptane. (D) 3,5-Bis(trichloromethyl)heptane.
3. The correct IUPAC name of the following compound is :



- (A) 5,6-Diethyl-8-methyldec-6-ene (B) 5,6-Diethyl-3-methyloct-4-ene
 (C) 5,6-Diethyl-3-methyldec-4-ene (D) 2,4,5-Triethylnon-3-ene
4. Diethenyl pentadiene is :



- 5.* Which IUPAC name is **correct** :

- (A) $\text{CH}_3-\underset{\text{C}_2\text{H}_5}{\text{CH}}=\underset{\text{C}_2\text{H}_5}{\text{CH}}-\text{CH}_3$ 2,3-Diethylbutene
- (B) $\text{HC}\equiv\text{C}-\underset{\text{HC}=\text{CH}_2}{\text{CH}}-\text{CH}=\text{CH}_2$ 3-Ethynylpenta-1, 4-diene
- (C) $\text{HC}\equiv\text{C}-\text{CH}=\text{CH}_2$ Butenyne
- (D) $\text{CH}_3-\text{CH}=\text{CH}-\text{C}\equiv\text{CH}$ Pent-3-en-1-yne

ORGANIC CHEMISTRY

DPP
DAILY PRACTICE PROBLEMS

DPP No. 5

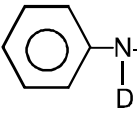
Total Marks : 31

Max. Time : 34 min.

Topic : IUPAC Nomenclature & Isomerism

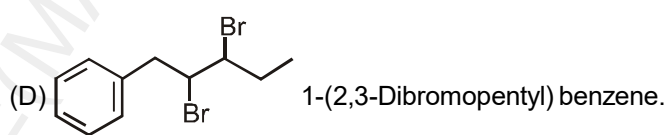
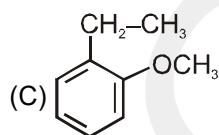
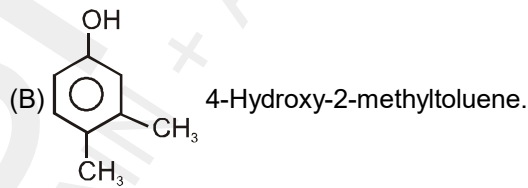
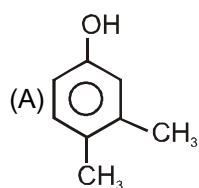
Type of Questions

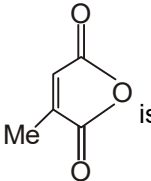
| Type of Questions | M.M., Min. |
|--|----------------------------|
| Single choice Objective ('-1' negative marking) Q.1 to Q.3 | (3 marks, 3 min.) [9, 9] |
| Assertion and Reason (no negative marking) Q.4 | (3 marks, 3 min.) [3, 3] |
| Multiple choice objective ('-1' negative marking) Q.5 | (4 marks, 4 min.) [4, 4] |
| Comprehension ('-1' negative marking) Q.6 | (3 marks, 3 min.) [3, 3] |
| Subjective Questions ('-1' negative marking) Q.7 | (4 marks, 5 min.) [4, 5] |
| Match the Following (no negative marking) Q.8 | (8 marks, 10 min.) [8, 10] |

1. IUPAC name of  is

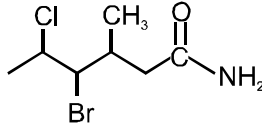
- (A) N-Deutero-N-formylbenzenamine (B) N-Phenylamino-N-deuteromethanal
(C) N-Deutero-N-phenylmethanamide (D) N-Deuterobenzene carboxamide

2. Which of the following is correct IUPAC name :



3. Correct IUPAC name of the compound  is

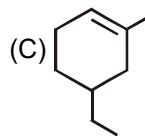
- (A) 2-Methylbutenedioic anhydride (B) 3-Methylbutenedioic anhydride
(C) 2-Methyl-1,4-diketobutene epoxy (D) 2-Methylcyclopentanoxy-1,4-dione

4. **Statement-1 :**  3-Bromo-2-chloro-4-methylpentanamide is incorrect IUPAC name.

Statement-2 : In case of chain terminating senior most functional group numbering starts from itself.

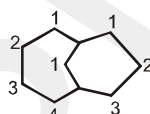
- (A) Statement-1 is True, Statement-2 is True; Statement-2 is a correct explanation for Statement-1.
(B) Statement-1 is True, Statement-2 is True; Statement-2 is NOT a correct explanation for Statement-1.
(C) Statement-1 is True, Statement-2 is False.
(D) Statement-1 is False, Statement-2 is True.

5.* Which of the following is/ are incorrect IUPAC name :

- (A) $\text{CH}_3 - \overset{\text{O}}{\parallel}{\text{C}} - \underset{\text{CH}_3}{\text{CH}} - \text{CH}_3$ 2-Methylbutan -3-one
- (B) $\text{CH}_3 - \underset{\text{OH}}{\text{CH}} - \text{COOH}$ 2-Hydroxypropanoic acid
- (C)  5-Ethyl-1-methylcyclohex-1-ene
- (D) $\text{CH}_3 - \underset{\text{CH}_3}{\text{CH}} - \overset{\text{O}}{\parallel}{\text{C}} - \underset{\text{CH}_3}{\text{CH}} - \text{OH}$ 4-Methyl-3-oxopentan-2-ol

6. **Comprehension #**

Bicyclo compounds are named by using the alkane name to designate the total number of carbon and bicyclo is used as prefix. While naming the bicycloalkane we write an expression between the word bicyclo and alkane (in square bracket), that denotes the number of carbon atoms in each bridge. The numerals are written in descending order and the numbers are separated by full stops.



Bicyclo [4. 3. 1] decane

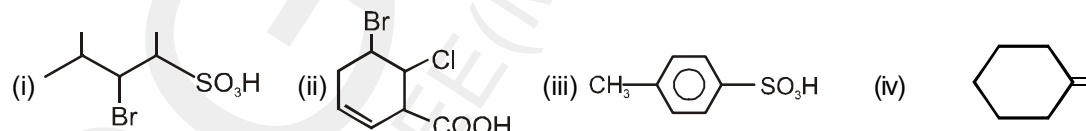
(a) The compound  is known by which of the following name :

- (A) Bicyclo [2. 2. 2] octane (B) Bicyclo [2. 2. 1] octane
 (C) Bicyclo [2. 2. 1] heptane (D) Bicyclo [2. 1. 2] heptane

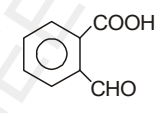
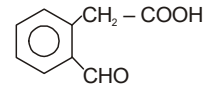
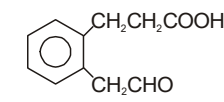
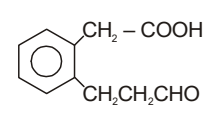
(b) The structure of bicyclo [1. 1. 0] butane is :

- (A)  (B)  (C)  (D) 

7. Write the IUPAC name of the following compound :



8. Match the following compound with their IUPAC name :

| Compound | IUPAC Name |
|---|--|
| (P)  | (W) 3-[2-(2-Oxoethyl)phenyl]propanoic acid |
| (Q)  | (X) 2-[2-(3-Oxopropyl)phenyl]ethanoic acid |
| (R)  | (Y) 2-(2-Formylphenyl)ethanoic acid |
| (S)  | (Z) 2-Formylbenzenecarboxylic acid |

ORGANIC CHEMISTRY

DPP
DAILY PRACTICE PROBLEMS

DPP No. 6

Total Marks : 32

Max. Time : 35 min.

Topic : IUPAC Nomenclature & Isomerism, Structural Isomerism

Type of Questions

Single choice Objective ('-1' negative marking) Q.1 to Q.4

(3 marks, 3 min.)

M.M., Min.

[12, 12]

Multiple choice objective ('-1' negative marking) Q.5 to Q.6

(4 marks, 4 min.)

[8, 8]

Subjective Questions ('-1' negative marking) Q.7

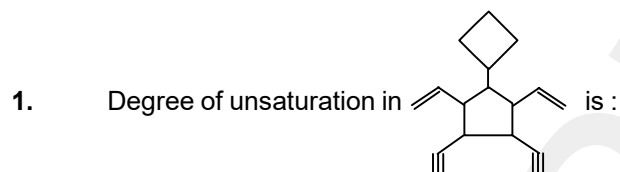
(4 marks, 5 min.)

[4, 5]

Match the Following (no negative marking) Q. 8

(8 marks, 10 min.)

[8, 10]



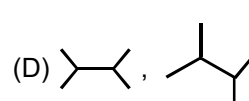
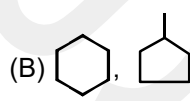
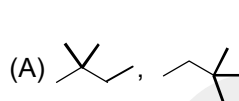
(A) 6

(B) 8

(C) 7

(D) 10

2. Which of the following pair is the chain isomer ?



3. Molecular formula C_4H_8 which type of isomerism will not show :

(A) Chain

(B) Metamerism

(C) Position

(D) Geometrical

4. Minimum carbon atoms are required to ketone to show position isomerism :

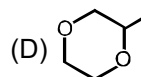
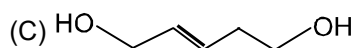
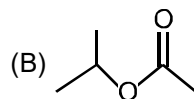
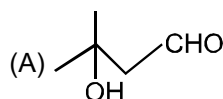
(A) 3

(B) 4

(C) 5

(D) 6

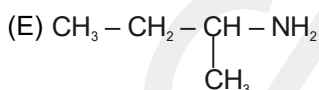
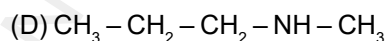
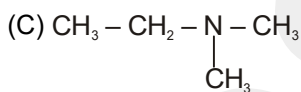
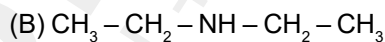
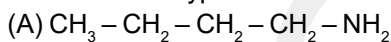
5.*  have functional isomer relation with



6.* Which is/are correctly matched.

| | | | | |
|-----|--|---|--|--------------------------|
| (A) | | , | | Metamers |
| (B) | | , | | Functional group isomers |
| (C) | | , | | Chain isomers |
| (D) | | , | | Positional isomers |

7. Fill the correct type of isomerism



(i) A & B are _____

(ii) B & C are _____

(iii) A & C are _____

(iv) B & D are _____

(v) A & E are _____

(vi) C & E are _____

8. Match the following:

| Column I Compound | Column II Degree of unsaturation |
|---|-------------------------------------|
| (A) | (p) 2 |
| (B) $\text{H}_3\text{C} - \text{CH}(\text{CH}_3) - \text{CH}_2 - \text{C}(\text{CH}_3) = \text{CH}_2$ | (q) 5 |
| (C) | (r) 4 |
| (D) | (s) 3 |

ORGANIC CHEMISTRY
DPP

DAILY PRACTICE PROBLEMS

DPP No. 7

Total Marks : 27

Max. Time : 28 min.

Topic : IUPAC Nomenclature & Isomerism
Type of Questions

| Type of Questions | M.M., Min. |
|--|----------------------------|
| Single choice Objective ('-1' negative marking) Q.1 to Q.5 | (3 marks, 3 min.) [15, 15] |
| Multiple choice objective ('-1' negative marking) Q.6 to Q.7 | (4 marks, 4 min.) [8, 8] |
| Subjective Questions ('-1' negative marking) Q.8 | (4 marks, 5 min.) [4, 5] |

- In allene (C_3H_4), the type(s) of hybridisation of the carbon atoms is (are) :
 (A) sp and sp^3 (B) sp and sp^2
 (C) only sp^3 (D) sp^2 and sp^3
- The carboxyl functional group ($-COOH$) is present in :
 (A) picric acid (B) barbituric acid
 (C) ascorbic acid (D) aspirin
- Aspirin is known as :
 (A) Acetyl salicylic acid (B) Phenyl salicylate
 (C) Acetyl salicylate (D) Methyl salicylic acid
- How many benzylic hydrogen are present in cumene :
 (A) 1 (B) 2 (C) 3 (D) 6
- How many allylic hydrogen are present in Methylbut-2-ene.
 (A) 6 (B) 8 (C) 9 (D) 10
- Which of the following is/are allylic alcohols ?
 (A) But-3-en-1-ol (B) But-3-en-2-ol
 (C) But-2-en-1-ol (D) But-1-en-1-ol
- Which name/names is/are correct ?
 (A) CH_2Cl_2 ; Methylene chloride (B) CH_3-CHCl_2 ; Ethylidene chloride
 (C) $CH_2 = CH - Cl$; Vinyl chloride (D) $HC \equiv C - CH_2 - Cl$; Propargyl chloride
- Write the structure of the following.
 (a) Benzoquinone (b) m-Xylene
 (c) Succinic acid (d) Lactic acid

ORGANIC CHEMISTRY

DPP

DAILY PRACTICE PROBLEMS

DPP No. 8

Total Marks : 30

Max. Time : 33 min.

Topic : Structural Determination

Type of Questions

Single choice Objective ('-1' negative marking) Q.1 to Q.5

(3 marks, 3 min.)

M.M., Min.

[15, 15]

Comprehension ('-1' negative marking) Q.6

(3 marks, 3 min.)

[3, 3]

Subjective Questions ('-1' negative marking) Q.7

(4 marks, 5 min.)

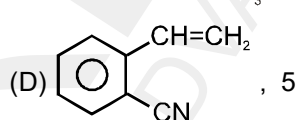
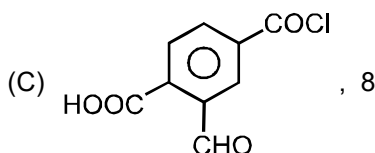
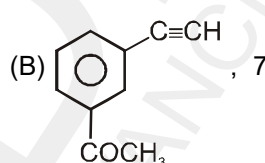
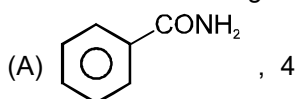
[4, 5]

Match the Following (no negative marking) Q.8

(8 marks, 10 min.)

[8, 10]

1. Which of the following is correctly matched with degree of unsaturation ?



2. How many alkene isomers will produce 1-Ethyl-3-methylcyclopentane on catalytic hydrogenation ?

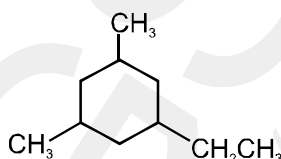
(A) 6

(B) 7

(C) 8

(D) 9

3. How many products (structural isomers) are formed by monochlorination of ?

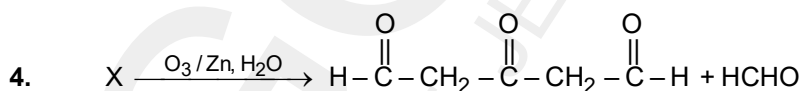


(A) 6

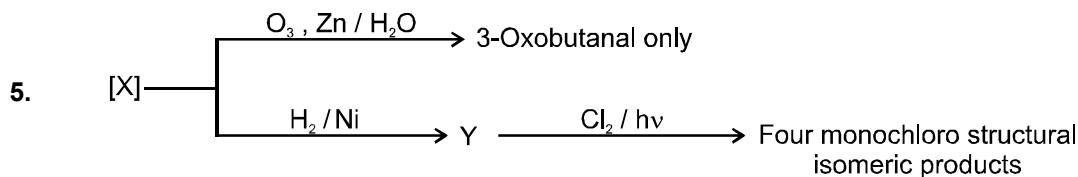
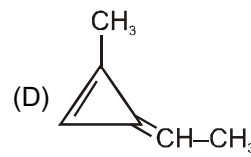
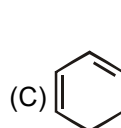
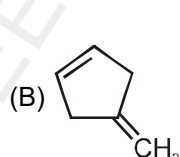
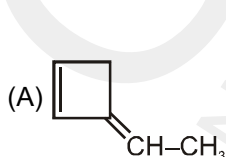
(B) 7

(C) 8

(D) 9



The structure of X will be :



Compound 'X' is :

(A) 1-Methylcyclopropene

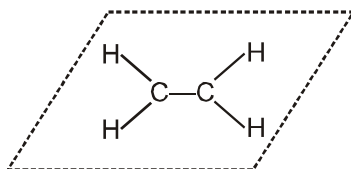
(B) 1, 4-Dimethylcyclohexa-1,4-diene

(C) 1, 4-Dimethylcyclohexa-1,3-diene

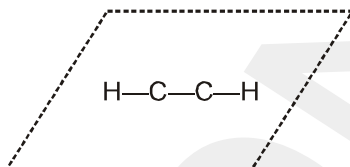
(D) 1, 2-Dimethylcyclohexa-1,4-diene

2. Draw the orbital diagrams for the following important structures.

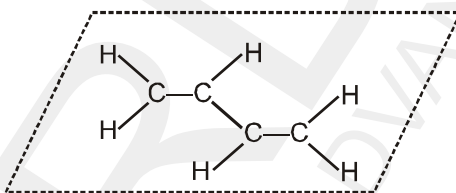
1. $\text{CH}_2 = \text{CH}_2$
 Ethyne



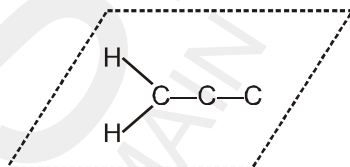
2. $\text{HC} \equiv \text{CH}$
 Ethyne



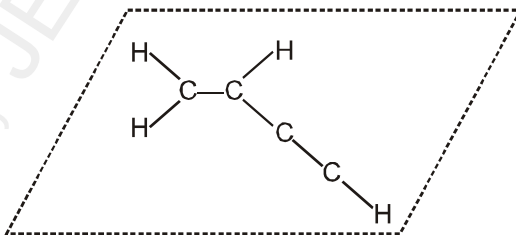
3. $\text{H}_2\text{C} = \text{CH} - \text{CH} = \text{CH}_2$
 Buta-1,3-diene



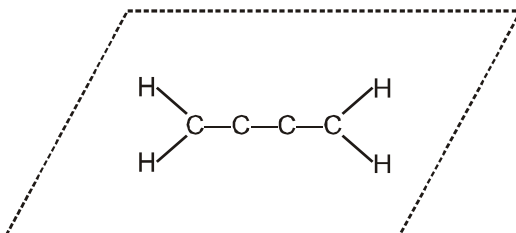
4. $\text{CH}_2 = \text{C} = \text{CH}_2$
 Propadiene



5. $\text{CH}_2 = \text{CH} - \text{C} \equiv \text{CH}$
 Butenyne



6. $\text{H}_2\text{C} = \text{C} = \text{C} = \text{CH}_2$
 Buta-1,2,3-triene



ORGANIC CHEMISTRY

DPP

DAILY PRACTICE PROBLEMS

DPP No. 10

Total Marks : 49

Max. Time : 50 min.

Topic : General Organic Chemistry

Type of Questions

Single choice Objective ('-1' negative marking) Q.1 to Q.15

(3 marks, 3 min.)

M.M., Min.

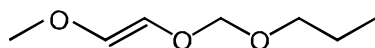
[45, 45]

Subjective Questions ('-1' negative marking) Q.16

(4 marks, 5 min.)

[4, 5]

1. How many lone pairs of electrons are there in the given compound ?



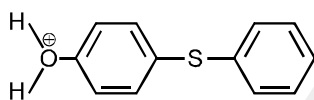
(A) 4

(B) 2

(C) 8

(D) 6

2. How many lone pairs of electrons are there in the given compound ?



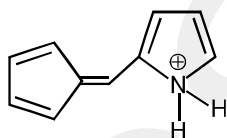
(A) 4

(B) 3

(C) 8

(D) 6

3. What is the hybridisation of positively charged nitrogen atom ?



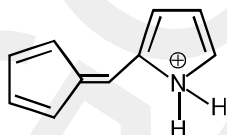
(A) sp

(B) sp²

(C) sp³

(D) None of these

4. How many sp² hybridised atoms are there in the given cation ?



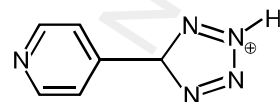
(A) 8

(B) 9

(C) 10

(D) 12

5. How many lone pairs are present in the given cation ?



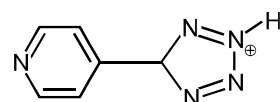
(A) 3

(B) 4

(C) 5

(D) 6

6. How many N atoms are sp² hybridised in the given cation ?



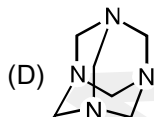
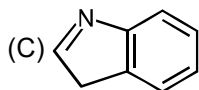
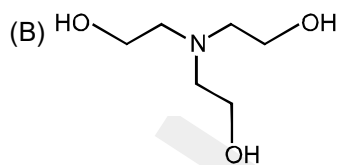
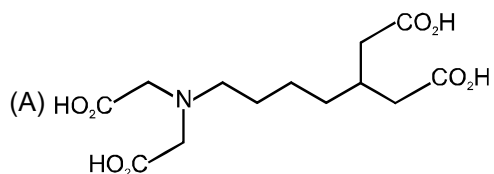
(A) 3

(B) 4

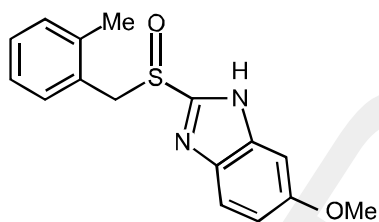
(C) 5

(D) 6

7. In which of the following molecules is the nitrogen atom sp^2 hybridised ?



8. Geometry around how many carbon atoms is tetrahedral in the given structure ?



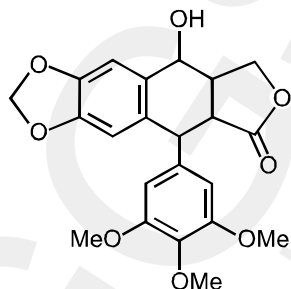
(A) 4

(B) 2

(C) 3

(D) 5

9. How many sp^3 hybridised carbon atoms are there in the given anti-cancer compound (podophyllotoxin) ?



the anti-cancer compound
podophyllotoxin

(A) 6

(B) 7

(C) 9

(D) 8

10. Bond order of C–C bond in benzene is :

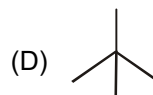
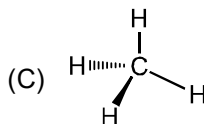
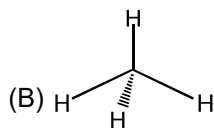
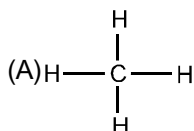
(A) 1

(B) 2

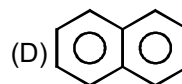
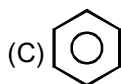
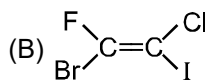
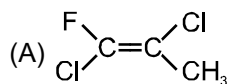
(C) 1.5

(D) two of above

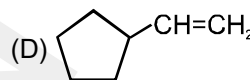
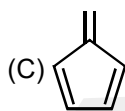
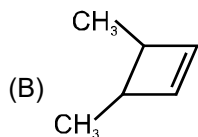
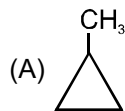
11. Which of the following is correct three dimensional representation of CH_4 ?



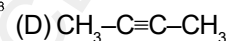
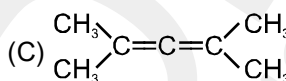
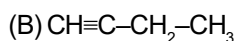
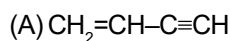
12. In which of the following all atoms do not present in the same plane ?



13. In which of the following all carbon atoms are present in the same plane ?



14. In which of the following are all C atoms linearly arranged ?



15. In How many carbon atoms are linearly arranged ?

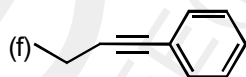
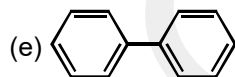
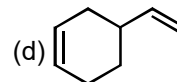
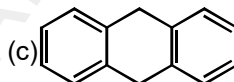
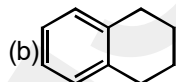
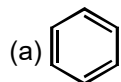
(A) 1

(B) 8

(C) 3

(D) 7

16. A maximum of how many (i) atoms (ii) carbon atoms must lie in same plane in each of the following molecule?



No. of $\text{sp}^2\text{-sp}^2$

σ -bonds

No. of $\text{sp}^2\text{-sp}$

σ -bonds

ORGANIC CHEMISTRY

DPP

DAILY PRACTICE PROBLEMS

DPP No. 11

Total Marks : 27

Max. Time : 29 min.

Topic : General Organic Chemistry

Type of Questions

Single choice Objective ('-1' negative marking) Q.1 to Q.5

(3 marks, 3 min.)

M.M., Min.

[15, 15]

Multiple choice objective ('-1' negative marking) Q.6

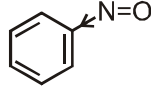
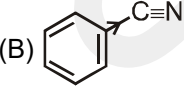
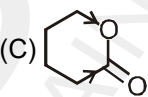
(4 marks, 4 min.)

[4, 4]

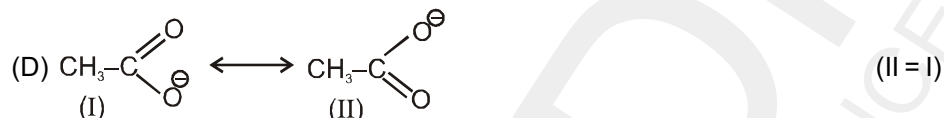
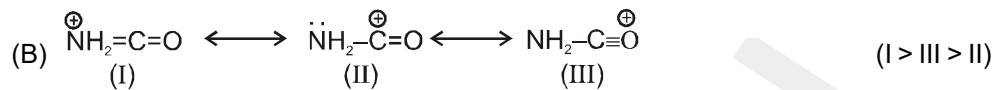
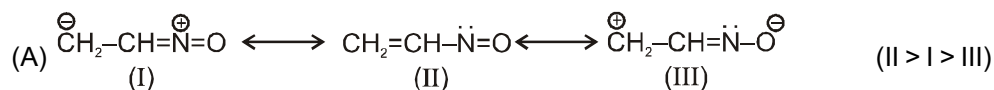
Subjective Questions ('-1' negative marking) Q.7 to Q.8

(4 marks, 5 min.)

[8, 10]

- Inductive effect is a permanent effect and is distance dependent.
 (A) Always (B) Some time (C) never (D) Can not decide
- Which of the following statement is CORRECT regarding the inductive effect ?
 (A) electron-donating inductive effect(+I effect) is generally more powerful than electron-withdrawing inductive effect(-I effect)
 (B) it implies the shifting of σ electrons from more electronegative atom to the lesser electronegative atom in a molecule
 (C) it implies the shifting of σ electrons from less electronegative atom to the more electronegative atom in a molecule
 (D) it increases with increase in distance.
- In which of the following species, incorrect direction of Inductive effect is/are shown ?
 (A)  (B)  (C)  (D) $\text{CH}_3 - \text{CH}_2 \leftarrow \text{MgBr}$
- Maximum -I effect is exerted by the group
 (A) C_6H_5 (B) $-\text{OCH}_3$ (C) $-\text{Cl}$ (D) $-\text{NH}_2$
- Which order of I effect is incorrect.
 (I) $-\overset{\oplus}{\text{N}}(\text{CH}_3)_3 > -\overset{\oplus}{\text{S}}(\text{CH}_3)_2$ [-I] (II) $-\text{OCH}_3 > -\text{OH}$ [-I]
 (III) $-\text{F} > -\text{Cl}$ [-I] (IV) $-\text{CH}_3 > -\overset{\ominus}{\text{O}}$ [+I]
 (A) II, III & IV (B) III & IV (C) IV only (D) all
- * Which of the following statement/s is/are correct for the inductive effect ?
 (A) It is a permanent effect (B) It transmits through sigma electrons
 (C) It is represented by \longleftrightarrow (D) It is represented by \longrightarrow or \longleftarrow .
- In which C - C bond of $\overset{3}{\text{C}}\text{H}_3 - \overset{2}{\text{C}}\text{H}_2 - \overset{1}{\text{C}}\text{H}_2 - \text{Br}$, the inductive effect is expected to be the least.
- How many groups show -I effect?
 $-\text{CH}_3$, $-\overset{\oplus}{\text{N}}\text{H}_3$, $-\text{OH}$, $-\text{O}^\ominus$, $-\text{N}(\text{CH}_3)_2$, $-\text{SO}_3\text{H}$, $-\text{CHO}$, $-\text{Cl}$, $-\text{COO}^\ominus$

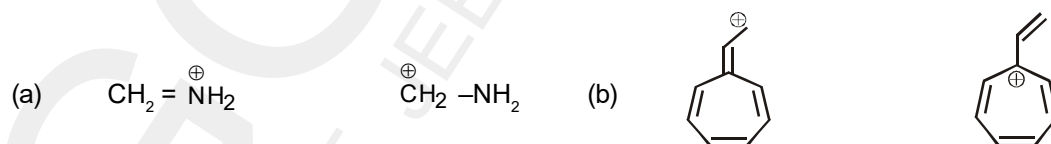
6.* Which of the following are correctly ordered for resonance stability



7. Match the resonance contributors in Column I with their attributes (properties) mentioned in Column II

| Column I | Column II |
|---|-----------------------|
| (A) $\text{CH}_3\text{CH}_2-\overset{\text{O}}{\parallel}{\text{C}}-\text{OCH}_2\text{CH}_3$ | (p) Equal contributor |
| (B) $\text{CH}_3-\overset{\oplus}{\text{N}}=\overset{\ominus}{\text{O}}$ | (q) major contributor |
| (C) $\text{CH}_3-\overset{\text{O}}{\parallel}{\text{C}}-\overset{\ominus}{\text{C}}\text{H}-\overset{\text{O}}{\parallel}{\text{C}}-\text{CH}_3$ | (r) minor contributor |

8. Give stability order in the following pairs.



ORGANIC CHEMISTRY

DPP
DAILY PRACTICE PROBLEMS

DPP No. 13

Total Marks : 30

Max. Time : 33 min.

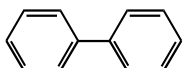
Topic : General Organic Chemistry

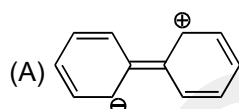
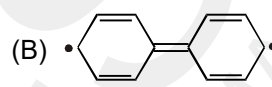
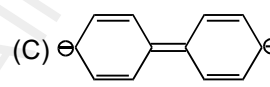
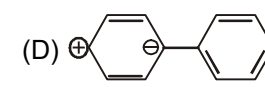
Type of Questions

| Type of Questions | M.M., Min. |
|--|----------------------------|
| Single choice Objective ('-1' negative marking) Q.1 to Q.5 | (3 marks, 3 min.) [15, 15] |
| Assertion and Reason (no negative marking) Q.6 | (3 marks, 3 min.) [3, 3] |
| Match the Following (no negative marking) Q.7 | (8 marks, 10 min.) [8, 10] |
| Subjective Questions ('-1' negative marking) Q.8 | (4 marks, 5 min.) [4, 5] |

1. Which of the following is/are resonating structures of diazomethane (CH_2N_2).

- (A) $\text{HN}=\text{C}=\text{NH}$ (B) $\text{CH}_2 = \overset{+}{\text{N}} = \text{N}^-$ (C) $\overset{-}{\text{C}}\text{H}_2 - \overset{+}{\text{N}} \equiv \text{N}$ (D) all of these

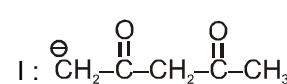
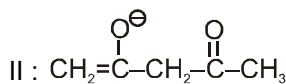
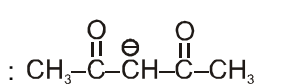
2. Which of the following does not represent the resonating structure of 

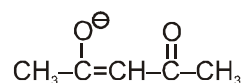
- (A)  (B)  (C)  (D) 

3. Decreasing +m-power of given group is :

- (I) $-\text{NH}_2$ (II) $-\text{OH}$ (III) $-\text{O}^-$ (IV) $-\text{NH}-\text{CO}-\text{CH}_3$
 (A) $\text{I} > \text{III} > \text{IV} > \text{II}$ (B) $\text{III} > \text{II} > \text{I} > \text{IV}$ (C) $\text{III} > \text{I} > \text{II} > \text{IV}$ (D) $\text{II} > \text{I} > \text{IV} > \text{III}$


4. The stability order of the following species is :

- I :  II :  III :  IV :



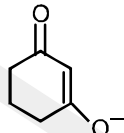
- (A) $\text{I} > \text{II} > \text{III} > \text{IV}$ (B) $\text{III} > \text{I} > \text{II} > \text{IV}$ (C) $\text{IV} > \text{II} > \text{III} > \text{I}$ (D) $\text{IV} > \text{III} > \text{II} > \text{I}$

5. Identify the correct statements

(i) All C – C bonds in  are equal.

(ii) All C – C bonds in $\text{CH}_2 = \text{CH} - \text{CH} = \text{CH}_2$ are equal.

(iii) All C – O bonds in $\text{CH}_3 - \text{C} \begin{matrix} \text{=O} \\ \text{O}^- \end{matrix}$ are equal.

(iv) All C – O bond in  are equal.

(A) i, ii, iii, iv

(B) i, iii, iv

(C) i, ii, iii

(D) ii, iii, iv

6. **STATEMENT -1** : Bond length of double bond in benzene is more than the bond length of double bond in buta-1,3-diene.

STATEMENT -2 : Increase in delocalisation of π electrons increases the bond length of double bond.

(A) Statement-1 is True, Statement-2 is True; Statement-2 is a correct explanation for Statement-1.

(B) Statement-1 is True, Statement-2 is True; Statement-2 is NOT a correct explanation for Statement-1

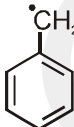
(C) Statement-1 is True, Statement-2 is False (D) Statement-1 is False, Statement-2 is True

7. Match the column :

Column-I

(A) $\text{CH}_2^+ - \text{CH} = \text{CH}_2$

(B) $\text{H}_2\text{N} - \text{CH} = \text{CH}_2$

(C) 

(D) $\text{H}_2\text{N} - \text{C}(\text{NH}) - \text{NH}_2$

Column-II

(p) Resonance possible

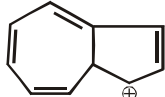
(q) Even number of p-electrons

(r) localized lone pair of e^- .

(s) Delocalized lone pair of e^- .

(t) $2 e^-$ in p orbitals

8. Find the total number of carbon where positive charge can be delocalised by true resonance [Including the given structure] :

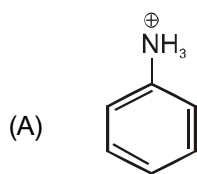
(a) 

(b) $\text{CH}_3 - \text{CH}^+ - \text{CH} = \text{CH} - \text{C}(\text{CH}_2) = \text{C}_6\text{H}_5$

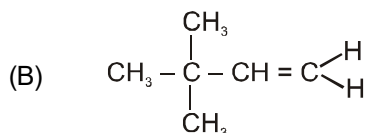
7. Match the compounds given in column I with their electronic effects mentioned in column II

Column I

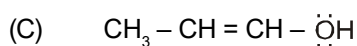
Column II



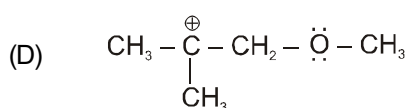
(p) Inductive effect



(q) Delocalisation of π electron

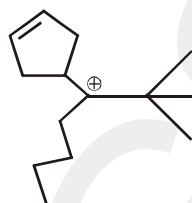


(r) Hyperconjugation



(s) Mesomeric effect

8. The total number of contributing structures showing hyperconjugation (involving C-H bonds) for the following molecule is



ORGANIC CHEMISTRY

DPP

DAILY PRACTICE PROBLEMS

DPP No. 15

Total Marks : 34

Max. Time : 40 min.

Topic : General Organic Chemistry

Type of Questions

Single choice Objective ('-1' negative marking) Q.1 to Q.2

(3 marks, 3 min.)

M.M., Min.

[6, 6]

Multiple choice objective ('-1' negative marking) Q.3

(4 marks, 4 min.)

[4, 4]

Match the Following (no negative marking) Q.4

(8 marks, 10 min.)

[8, 10]

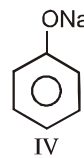
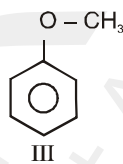
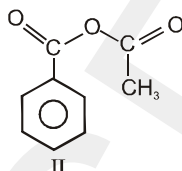
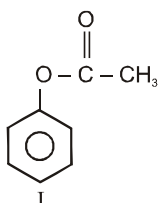
Subjective Questions ('-1' negative marking) Q.5 to Q.8

(4 marks, 5 min.)

[16, 20]

1. The reactivity order of benzene ring for the given reaction is (benzene ring with highest π electron density will

be most reactive) Aromatic compound $\xrightarrow{E^{\oplus}}$



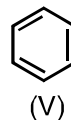
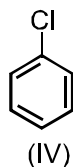
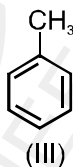
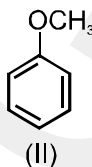
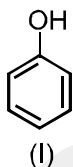
(A) III > I > IV > II

(C) III > IV > II > I

(B) I > III > IV > II

(D) IV > III > I > II

2. Arrange the following compounds in the order of decreasing reactivity towards electrophilic substitution



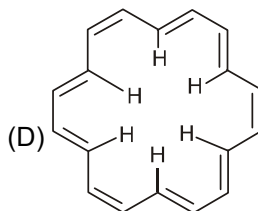
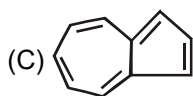
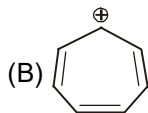
(A) V > IV > III > II > I

(C) I > II > IV > III > V

(B) I > II > III > V > IV

(D) I > III > IV > II > V

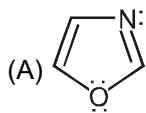
3.* Which of the following are Aromatic in nature.



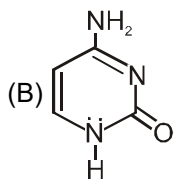
4. Match the following :

Column – I

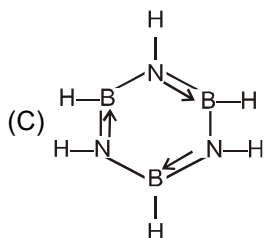
Column – II



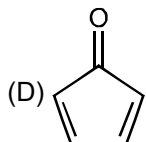
(p) Aromatic



(q) Non aromatic

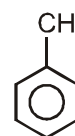
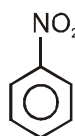
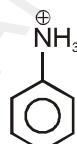
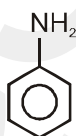
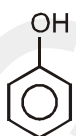
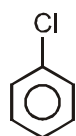


(r) Anti aromatic

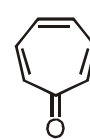
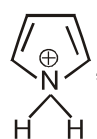
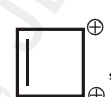


(s) Heterocyclic

5. How many of the following compounds are more reactive than benzene towards electrophilic substitution.

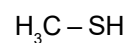
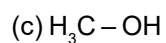
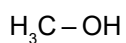
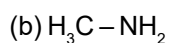
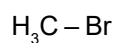
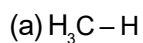



6. How many species out of the following are aromatic ?



7. Explain the terms Inductive and Electromeric effects.

8. Which bond is more polar in the following pairs of molecules?



| S.No. | Species | + I | - I | + m | - m | + H.C. | - H.C. |
|-------|---|-----|-----|-----|-----|--------|--------|
| 20 | -I | | | | | | |
| 21 | $-\overset{\oplus}{\text{N}}=\text{N}:$ | | | | | | |
| 22 | $-\text{C}\equiv\overset{\oplus}{\text{O}}:$ | | | | | | |
| 23 | -S-R | | | | | | |
| 24 | $\begin{array}{c} \text{O} \\ \\ -\text{S}-\text{R} \\ \\ \text{O} \end{array}$ | | | | | | |
| 25 | $-\overset{\oplus}{\text{P}}\text{R}_3$ | | | | | | |
| 26 | $-\overset{\oplus}{\text{N}}\text{R}_3$ | | | | | | |
| 27 | $-\text{CH}_3$ | | | | | | |
| 28 | $-\overset{\oplus}{\text{S}}\text{R}_2$ | | | | | | |
| 29 | $-\text{CMe}_3$ | | | | | | |
| 30 | $-\text{CF}_3$ | | | | | | |
| 31 | $-\text{CCl}_3$ | | | | | | |
| 32 | -Ph | | | | | | |
| 33 | $-\text{CH}=\text{CH}_2$ | | | | | | |
| 34 | $-\text{COOH}$ | | | | | | |
| 35 | $-\text{O}-\text{CH}_3$ | | | | | | |
| 36 | $-\ddot{\text{O}}\text{H}$ | | | | | | |
| 37 | $-\dot{\text{N}}\text{H}_2$ | | | | | | |
| 38 |  | | | | | | |
| 39 | $-\overset{\ominus}{\text{C}}\text{R}_2$ | | | | | | |
| 40 | $-\text{CR}_2$ | | | | | | |

ANSWER KEY (ORGANIC CHEMISTRY)

ORGANIC CHEMISTRY

DPP No. # 1

1. (C) 2. (C) 3. (C) 4. (D) 5*. (BCD) 6*.
 (ABD) 7. 9 8. (A → p, s) ; (B → p, q) ; (C → r, s) ; (D → p)

DPP No. # 2

1. (B) 2. (C) 3. (D) 4. (D) 5. (B)
 6. (A,B) 7. 114. 8. A → S, B → P, C → Q, D → R

DPP No. # 3

1. (C) 2. (C) 3. (A) 4. (D)
 5.* (A,B,D) 6.* (A,B,C,D)
 7. $H_2C = C = CH_2$
 M. W. = 40. 8. [A - s] ; [B - q] ; [C - p] ; [D - r]

DPP No. # 4

1. (A) 2. (D) 3. (C) 4. (B)
 5. (B,C,D) 6. (B,C,D) 7. 6
 8. (a) (B) (b) (C) (c) (D)

DPP No. # 5

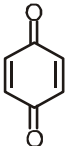
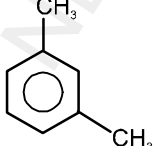
1. (C) 2. (A) 3. (A) 4. (A) 5.* (A, D) 6.
 (a) (C) (b) (C) 8. (P → Z) ; (Q → Y) ; (R → W) ; (S → X).

DPP No. # 6

1. (B) 2. (B) 3. (B) 4. (C)
 5. (ABCD) 6. (ABC) 8. (A → s) ; (B → r) ; (C → q) ; (D → p)

DPP No. # 7

1. (B) 2. (D) 3. (A) 4. (A) 5. (C)
 6. (B C) 7. (ABCD)

8. (a)  (b)  (c) $\begin{array}{c} CH_2 - COOH \\ | \\ CH_2 - COOH \end{array}$ (d) $\begin{array}{c} CH_3 - CH - COOH \\ | \\ OH \end{array}$

DPP No. # 8

1. (B) 2. (C) 3. (B) 4. (B) 5. (D)
 6. (a). (C) (b). (A) 8. (A → q) ; (B → s) ; (C → p) ; (D → r)

DPP No. # 9

Hyb. : $sp^3 \quad sp^3 \quad sp^3$
1. $CH_3-CH_2-CH_3$
 B. A $109^\circ \quad 109^\circ \quad 109^\circ$

Hyb. $sp^3 \quad sp^2 \quad sp^2$
2. $CH_3-CH=CH_2$
 B. A $109^\circ \quad 120^\circ \quad 120^\circ$

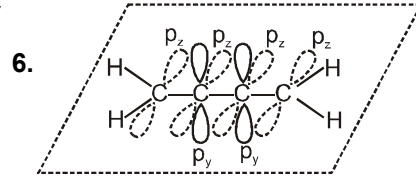
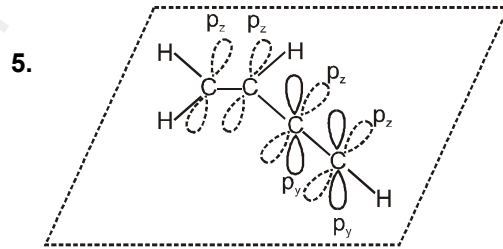
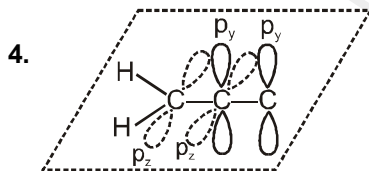
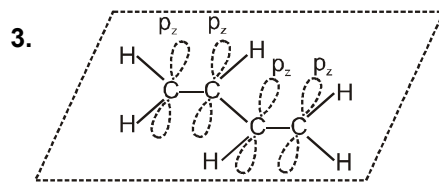
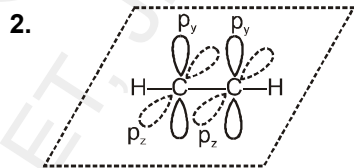
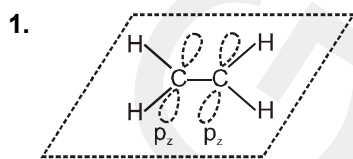
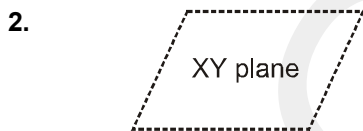
Hyb.: $sp^3 \quad sp \quad sp$
3. $CH_3-C \equiv CH$
 B. A $109^\circ \quad 180^\circ \quad 180^\circ$

Hyb. : $sp^2 \quad sp^2 \quad sp^2 \quad sp^2$
4. $CH_2=CH-CH=CH_2$
 B. A $120^\circ \quad 120^\circ \quad 120^\circ \quad 120^\circ$

Hyb. : $sp^2 \quad sp^2 \quad sp \quad sp$
5. $CH_2=CH-C \equiv CH$
 B. A $120^\circ \quad 120^\circ \quad 180^\circ \quad 180^\circ$

Hyb. : $sp^2 \quad sp \quad sp^2$
6. $CH_2=C=CH_2$
 B. A $120^\circ \quad 180^\circ \quad 120^\circ$

Hyb. : $sp^2 \quad sp^2 \quad sp^3 \quad sp^2 \quad sp \quad sp^2$
7. $CH_2=CH-CH_2-HC=C=CH_2$
 B. A $120^\circ \quad 120^\circ \quad 109^\circ \quad 120^\circ \quad 180^\circ \quad 120^\circ$



DPP No. # 10

- | | | | | |
|---------------|---------------|---------------|---------------|----------------|
| 1. (D) | 2. (B) | 3. (C) | 4. (C) | 5. (B) |
| 6. (C) | 7. (C) | 8. (C) | 9. (C) | 10. (C) |

11. (C) 12. (A) 13. (C) 14. (D) 15. (C)
16. (a) (i) 12, (ii) 6. (b) (i) 12, (ii) 8. (c) (i) 12, (ii) 8. (d). (i) 6, (ii) 4.
 (e) (i) 14, (ii) 8. (f) (i) 14, (ii) 9.

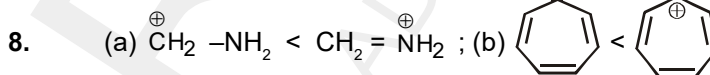
DPP No. # 11

1. (A) 2. (C) 3. (A) 4. (C) 5. (C)
- 6.* (ABD)
7. Inductive effect is expected to be the least in the bond between carbon 3 and carbon 2. 8. 6

DPP No. # 12

- 1.* (ABCD) 2.* (ABC) 3.* (ACD) 4. (B) 5. (D)
- 6.* (ABD)

7. (A - q) ; (B - p) ; (C - r)



DPP No. # 13

1. (B) 2. (B) 3. (C) 4. (D) 5. (B)
6. (A) 7. (A) - p,q,t ; (B) - p,q,s (C) - p (D) - p,q,r,s 8. (a) 5 (b) 3

DPP No. # 14

- 1.* (ABC) 2. (A) 3.* (A,D) 4. (A) 5. (C)
6. (C) 7. (A → p, q), (B → p), (C → p, q, r, s), (D → p, r) 8. 7

DPP No. # 15

1. (D) 2. (B) 3. (BCD)
4. (A - p,s) ; (B - p,s) ; (C - p,s) ; (D - r) 5. 3 6. 2



GGSRDN

Educational Services Private Limited

9th, 10th, NEET, JEE (Main/Advanced)

अभ्यास ही सबसे बड़ा गुरु है।

CLASS : XI

(ORGANIC CHEMISTRY)

DPP

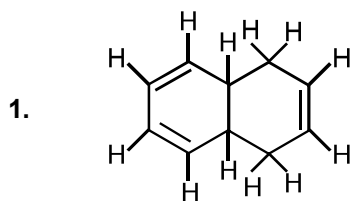
DAILY PRACTICE PROBLEM

Solutions

DPP-1 to 16

ORGANIC CHEMISTRY

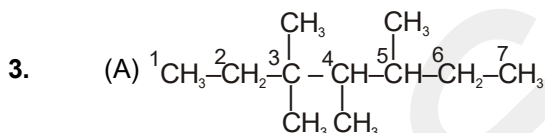
DPP No. # 1



DPP No. # 2

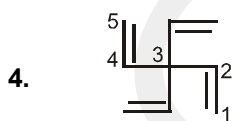
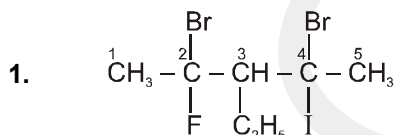
8. (A) $\text{CH}_3\text{-CH}_2\text{-CH}_2\text{-CH}_3$
 M.F. = $\text{C}_4\text{H}_{10} = 48 + 10 = 58$
 (B) $\text{HC}^\ominus\text{CH}$
 M.F. @ $\text{C}_2\text{H}_2 = 24 + 2 = 26$
 (C) $\text{CH}_3\text{-CH}=\text{CH}_2$
 M.F. @ $\text{C}_3\text{H}_6 = 36 + 6 = 42$
 (D) $\text{CH}_3\text{-C}^\ominus\text{CH}$
 M.F. @ $\text{C}_3\text{H}_4 = 36 + 4 = 40$

DPP No. # 3

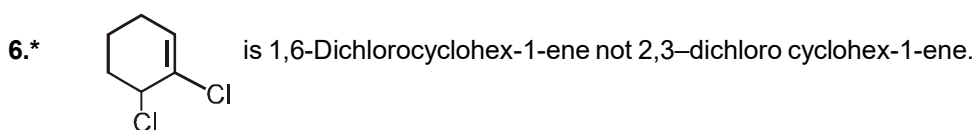
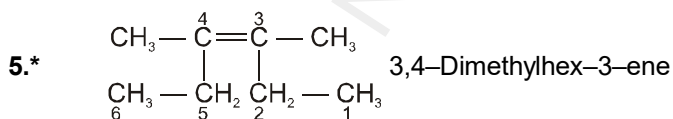


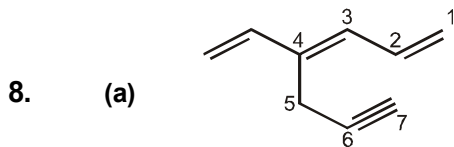
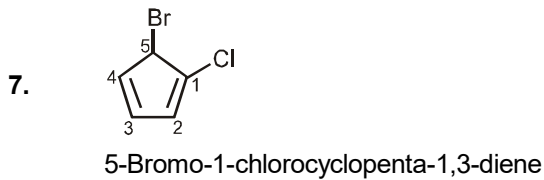
Lowest set of Locant (3, 3, 4, 5)

DPP No. # 4

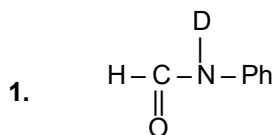


Diethenyl pentadiene

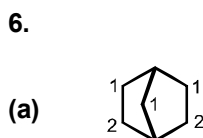




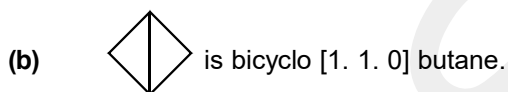
DPP No. # 5



N-Deutero-N-phenylmethanamide.



Bicyclo [2. 2. 1] heptane

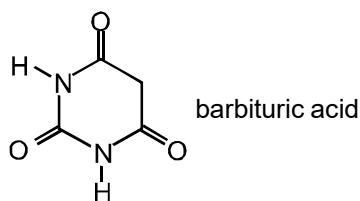
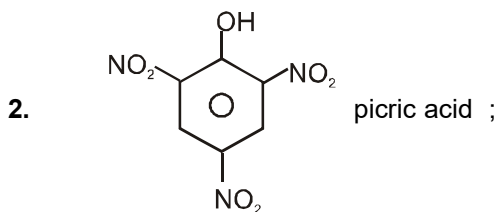
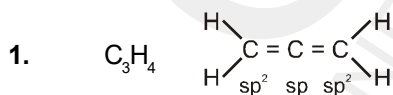


7. (i) 3-Bromo-4-methylpentane-2-sulphonic acid
 (ii) 5-Bromo-6-chlorocyclohex-2-ene-1-carboxylic acid
 (iii) 4-Methylbenzene-1-sulphonic acid
 (iv) Methylene cyclohexane

DPP No. # 6

7. (i) Functional isomer (ii) Functional isomer (iii) Functional isomer
 (iv) Metamer (v) Position isomer (vi) Functional isomer

DPP No. # 7



1.

Hyb. : $sp^3 \quad sp^3 \quad sp^3$

1. $CH_3-CH_2-CH_3$
 B. A $109^\circ \quad 109^\circ \quad 109^\circ$

Hyb. $sp^3 \quad sp^2 \quad sp^2$

2. $CH_3-CH=CH_2$
 B. A $109^\circ \quad 120^\circ \quad 120^\circ$

Hyb.: $sp^3 \quad sp \quad sp$

3. $CH_3-C \equiv CH$
 B. A $109^\circ \quad 180^\circ \quad 180^\circ$

Hyb. : $sp^2 \quad sp^2 \quad sp^2 \quad sp^2$

4. $CH_2=CH-CH=CH_2$
 B. A $120^\circ \quad 120^\circ \quad 120^\circ \quad 120^\circ$

Hyb. : $sp^2 \quad sp^2 \quad sp \quad sp$

5. $CH_2=CH-C \equiv CH$
 B. A $120^\circ \quad 120^\circ \quad 180^\circ \quad 180^\circ$

Hyb. : $sp^2 \quad sp \quad sp^2$

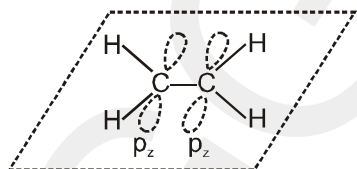
6. $CH_2=C=CH_2$
 B. A $120^\circ \quad 180^\circ \quad 120^\circ$

Hyb. : $sp^2 \quad sp^2 \quad sp^3 \quad sp^2 \quad sp \quad sp^2$

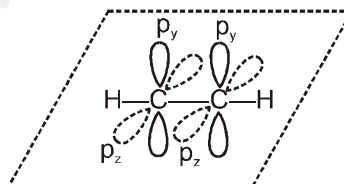
7. $CH_2=CH-CH_2-HC=C=CH_2$
 B. A $120^\circ \quad 120^\circ \quad 109^\circ \quad 120^\circ \quad 180^\circ \quad 120^\circ$

2.

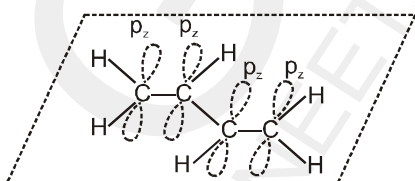
1.



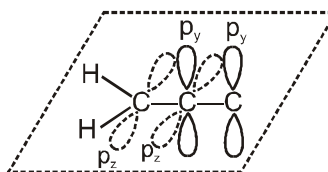
2.



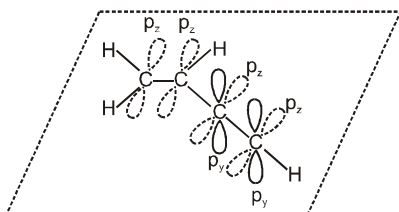
3.



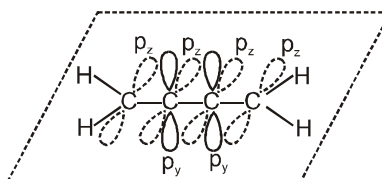
4.



5.



6.



DPP No. # 11

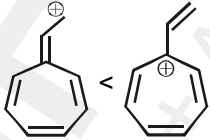
2. ex. $\overset{\delta+}{CH_3} \rightarrow \overset{\delta+}{CH_2} \rightarrow \overset{\delta-}{Cl}$

3. Case A has incorrect direction of I-effect.

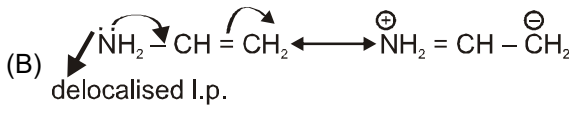
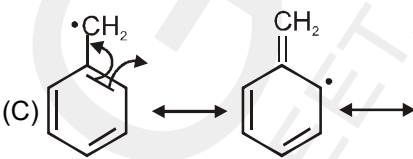
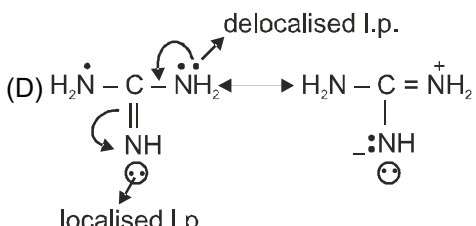
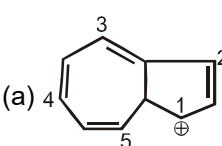
4. Maximum - I effect - Cl.

- Self explanatory.
- Self explanatory.
- Magnitude of inductive effect diminishes as the number of intervening bonds increases. Hence, the effect is least in the bond between carbon 3 and carbon 2.

DPP No. # 12

- (A,B,C,D)
Due to delocalization of paired & unpaired p-orbital electrons
- (A,B,C,) $\text{CH}_2 = \text{C} = \text{C} = \text{CH}_2 \longleftrightarrow \overset{\oplus}{\text{CH}_2} - \text{C} \equiv \text{C} - \overset{\ominus}{\text{CH}_2}$
 $\longleftrightarrow \overset{\ominus}{\text{CH}_2} - \text{C} \equiv \text{C} - \overset{\oplus}{\text{CH}_2}$
- Lone pair of electrons of $\text{H}_2\text{C} = \dot{\text{N}} - \text{CH}_3$ is in sp^2 hybrid orbital.**
 $\text{H}_2\text{C} = \dot{\text{N}} - \text{CH}_3$ dk ,dkdh bysDV^akWu ;qXe sp^2 ladfj r d{k d esa gSA
- * (C) $\text{CH}_3 - \overset{\oplus}{\text{C}} = \text{O} \longleftrightarrow \text{CH}_3 - \overset{\oplus}{\text{C}} \equiv \text{O}$
 (I) (II)
 Stability II > I due to No. of π bonds.
- (A - q) ; (B - p) ; (C - r)
- (a) $\overset{\oplus}{\text{CH}_2} - \text{NH}_2 < \text{CH}_2 = \overset{\oplus}{\text{NH}_2}$; (b) 

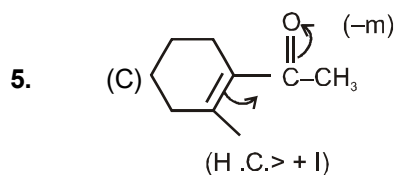
DPP No. # 13

- Self explanatory
- (Moderate)** There are unpaired electrons, others have no unpaired electrons.
- $-\overset{\ominus}{\text{O}} > -\text{NH}_2 > -\text{OH} > -\text{NH} - \text{CO} - \text{CH}_3$
- Due to delocalization of π electron in benzene.
- (A) $\overset{\oplus}{\text{CH}_2} - \text{CH} = \text{CH}_2 \longleftrightarrow \text{CH}_2 - \text{CH} = \overset{\oplus}{\text{CH}_2}$ (B) 
 delocalised l.p.
- (C) 
 (D) 
 delocalised l.p.
 localised l.p.
8. (a) 
 (b) $\text{CH}_3 - \underset{1}{\text{CH}} - \underset{2}{\text{CH}} = \underset{3}{\text{CH}} - \text{C}(\text{CH}_2) = \text{C}_6\text{H}_5$

DPP No. # 14

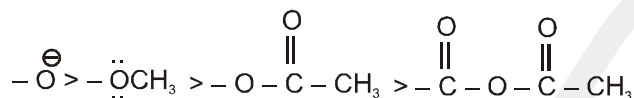
- (A,D) $-\text{CH}_3$ (H.C. & + I)
 $-\text{OCH}_3$ (+M > - I)

4. $\text{CH}_3 - \text{CH} = \text{CH} - \text{CH} = \text{CH} - \text{CH}_3$
 (A) reso + H.C.
 $\text{>C} = \text{C} <$ bond order \downarrow as, bond length $\uparrow \propto \sigma$

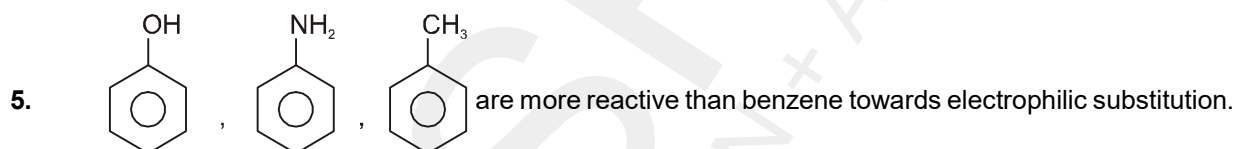


DPP No. # 15

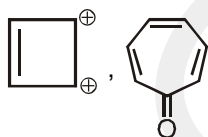
1. (D) Due to +m order



2. Rate of electrophilic substitution reaction \propto Stability of arenium ion.
 4. Aromatic \rightarrow planar, cyclic, $(4n+2) \pi e^-$, complete conjugation
 Antiaromatic \rightarrow planar, cyclic, $(4n) \pi e^-$, complete conjugation
 Non aromatic \rightarrow cyclic structure with non-planar geometry with any hybridization



6. Aromatic species are



7. Inductive effect is permanent displacement of shared pair of electron along the chain of carbon atom due to presence of polar covalent bond.
 Electromeric effect is a temporary effect. It is defined as the complete transfer of a shared pair of π -electrons to one of the atoms joined by a multiple bond on the demand of an attacking reagent.
8. (a) $\text{H}_3\text{C}-\text{Br}$ (b) $\text{H}_3\text{C}-\text{OH}$ (c) $\text{H}_3\text{C}-\text{OH}$

DPP No. # 16

| | | | | | | | |
|----|---|---|---|---|---|---|---|
| 23 | $-S-R$ | | ✓ | ✓ | ✓ | | |
| 24 | $\begin{array}{c} O \\ \\ -S-R \\ \\ O \end{array}$ | | ✓ | | ✓ | | |
| 25 | $^{\oplus}PR_3$ | | ✓ | | ✓ | | |
| 26 | $^{\oplus}NR_3$ | | ✓ | | | | |
| 27 | $-CH_3$ | ✓ | | | | ✓ | |
| 28 | $^{\oplus}SR_2$ | | ✓ | | ✓ | | |
| 29 | $-CMe_3$ | ✓ | | | | | |
| 30 | $-CF_3$ | | ✓ | | | | ✓ |
| 31 | $-CCl_3$ | | ✓ | | | | ✓ |
| 32 | $-Ph$ | | ✓ | ✓ | ✓ | | |
| 33 | $-CH=CH_2$ | | ✓ | ✓ | ✓ | | |
| 34 | $-COOH$ | | ✓ | | ✓ | | |
| 35 | $-O-CH_3$ | | ✓ | ✓ | | | |
| 36 | $-\ddot{O}H$ | | ✓ | ✓ | | | |
| 37 | $-\ddot{N}H_2$ | | ✓ | ✓ | | | |
| 38 | $-N$ (in a six-membered ring) | | ✓ | ✓ | | | |
| 39 | $^{\ominus}CR_2$ | ✓ | | ✓ | | | |
| 40 | $^{\oplus}CR_2$ | | ✓ | | ✓ | | |