

# IITJEE 2009

## Paper 1 Code (0)

### Chemistry

1. Given that the abundances of isotopes  $^{54}\text{Fe}$ ,  $^{56}\text{Fe}$  and  $^{57}\text{Fe}$  are 5%, 90% and 5%, respectively, the atomic mass of Fe is
- (A) 55.85      (B) 55.95      (C) 55.75      (D) 56.05

1: Atomic mass =  $\sum X_i \cdot A_i$

$$= 54 \times 0.05 + 56 \times 0.9 + 57 \times 0.05$$

$$= 55.95$$

Ans: (B)

2. The term that corrects for the attractive forces present in a real gas in the van der Waals equation is
- (A)  $nb$       (B)  $\frac{an^2}{V^2}$       (C)  $-\frac{an^2}{V^2}$       (D)  $-nb$

2: The term for the attractive forces present in a real gas in Van der Waals equation is:

$$\left( P + \frac{an^2}{V^2} \right) (V - nb) = nRT$$

$\frac{an^2}{V^2}$

Ans: (B)

3. Among the electrolytes  $\text{Na}_2\text{SO}_4$ ,  $\text{CaCl}_2$ ,  $\text{Al}_2(\text{SO}_4)_3$  and  $\text{NH}_4\text{Cl}$ , the most effective coagulating agent for  $\text{Sb}_2\text{S}_3$  sol is
- (A)  $\text{Na}_2\text{SO}_4$       (B)  $\text{CaCl}_2$       (C)  $\text{Al}_2(\text{SO}_4)_3$       (D)  $\text{NH}_4\text{Cl}$

3:  $Sb_2S_3$  is a -ively charged sol. To coagulate it most effectively, charge on cation should be maximum. Thus,  $Al^{3+}$  in  $Al_2(SO_4)_3$  will be most effective.

Ans : (c)

$1.0 \times 10^{-3}$  atm. The mole fraction of air dissolved in water at 298 K and 5 atm pressure is

- (A)  $4.0 \times 10^{-4}$       (B)  $4.0 \times 10^{-5}$       (C)  $5.0 \times 10^{-4}$       (D)  $4.0 \times 10^{-6}$

#### 4: Henry's law:

$$\rho_{\text{Partial}} = \rho_{\text{H}_2} \cdot (X_{N_2})_{\text{in Soln}}$$

$$5 \times 0.8 = 10 \times X_{N_2}$$

$P_{\text{Total}}$  ( $X_{N_2}$ ) in air

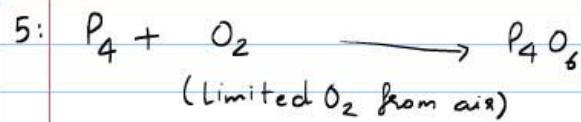
$$\Rightarrow X_{N_2} = 4 \times 10^{-5} = \frac{n_{N_2}}{n_{N_2} + n_{O_2}}$$

$$\approx \frac{n_{n_2}}{n_{\theta}}$$

$$\Rightarrow n_{N_2} = 4 \times 10^{-4}$$

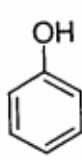
Ans : (A)

5 The reaction of  $P_4$  with  $X$  leads selectively to  $P_4O_6$ . The  $X$  is

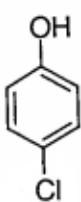


**Ans: (B)**

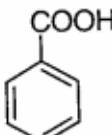
6. The correct acidity order of the following is



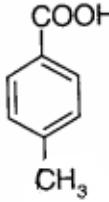
(I)



(II)

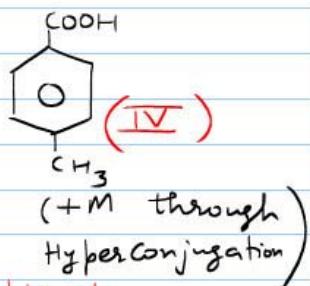
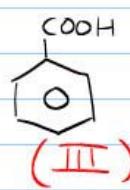
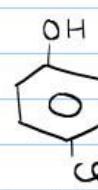
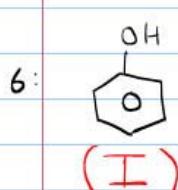


(III)



(IV)

- (A) (III) > (IV) > (II) > (I)      (B) (IV) > (III) > (I) > (II)  
 (C) (III) > (II) > (I) > (IV)      (D) (II) > (III) > (IV) > (I)



(-I, +M)

≈ Weak deactivating

(+M through)  
Hyperconjugation

\* -COOH is more acidic than phenols.

⇒ III > IV > II > I

**Ans: (A)**

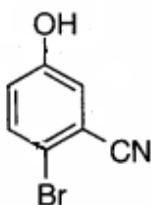
7. Among cellulose, poly(vinyl chloride), nylon and natural rubber, the polymer in which the intermolecular force of attraction is weakest is

- (A) Nylon      (B) Poly(vinyl chloride)  
 (C) Cellulose      (D) Natural Rubber

7: Natural rubber is a polymer of isoprene.  
Nylon, cellulose & PVC contains functional groups.

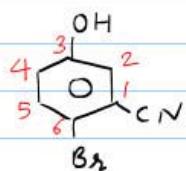
**Ans: (D)**

8. The IUPAC name of the following compound is



- (A) 4-Bromo-3-cyanophenol
- (B) 2-Bromo-5-hydroxybenzonitrile
- (C) 2-Cyano-4-hydroxybromobenzene
- (D) 6-Bromo-3-hydroxybenzonitrile

8:



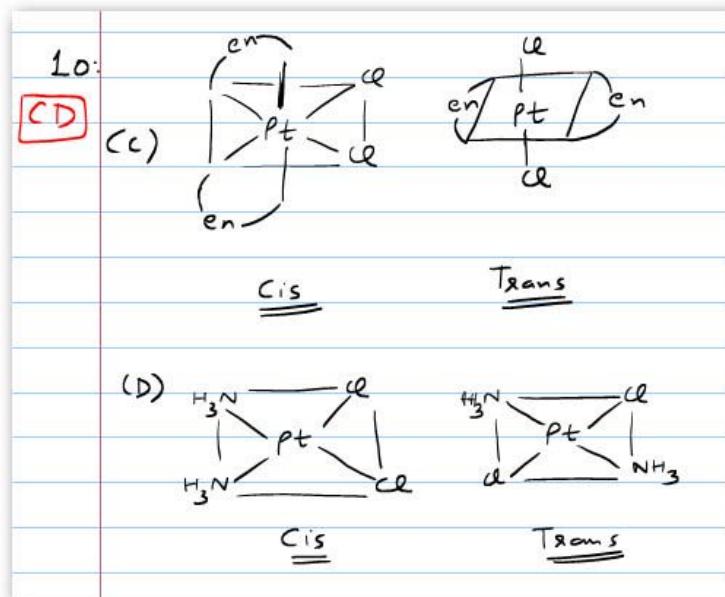
Clearly, the IUPAC name is:

6 - Bromo - 3 - hydroxybenzonitrile

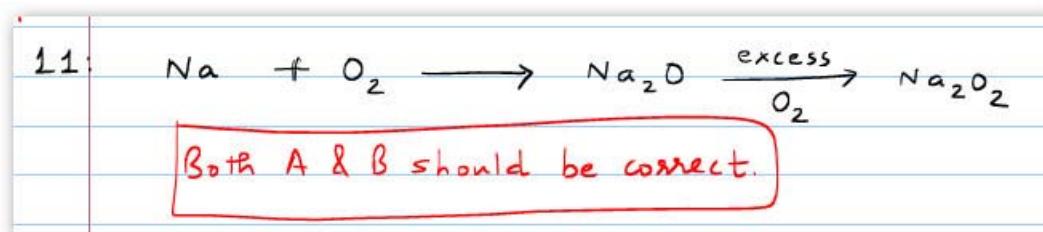
**Ans: (D)**

9. The correct statement(s) regarding defects in solids is(are)

- (A) Frenkel defect is usually favoured by a very small difference in the sizes of cation and anion
- (B) Frenkel defect is a dislocation defect
- (C) Trapping of an electron in the lattice leads to the formation of F-center
- (D) Schottky defects have no effect on the physical properties of solids

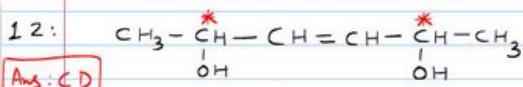


11. The compound(s) formed upon combustion of sodium metal in excess air is(are)  
(A)  $\text{Na}_2\text{O}_2$       (B)  $\text{Na}_2\text{O}$       (C)  $\text{NaO}_2$       (D)  $\text{NaOH}$



12. The correct statement(s) about the compound  $\text{H}_3\text{C}(\text{HO})\text{HC}-\text{CH}=\text{CH}-\text{CH}(\text{OH})\text{CH}_3$  (**X**) is(are)

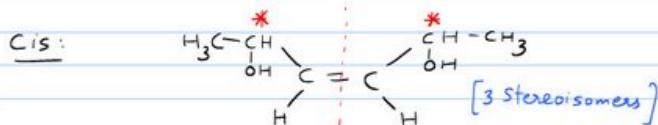
  - (A) The total number of stereoisomers possible for **X** is 6
  - (B) The total number of diastereomers possible for **X** is 3
  - (C) If the stereochemistry about the double bond in **X** is *trans*, the number of enantiomers possible for **X** is 4
  - (D) If the stereochemistry about the double bond in **X** is *cis*, the number of enantiomers possible for **X** is 2



Ans: CD

(x)

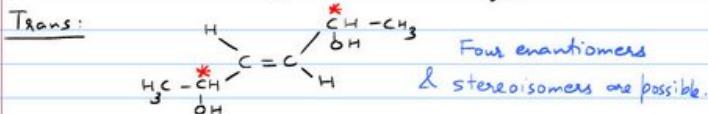
Total number of chiral centres = 2



↙ (Plane of symmetry)

(meso compound)

⇒ There are only two enantiomers for X.



⇒ Total stereoisomers = 3 + 4 = 7

There will be lot of diastereomers.

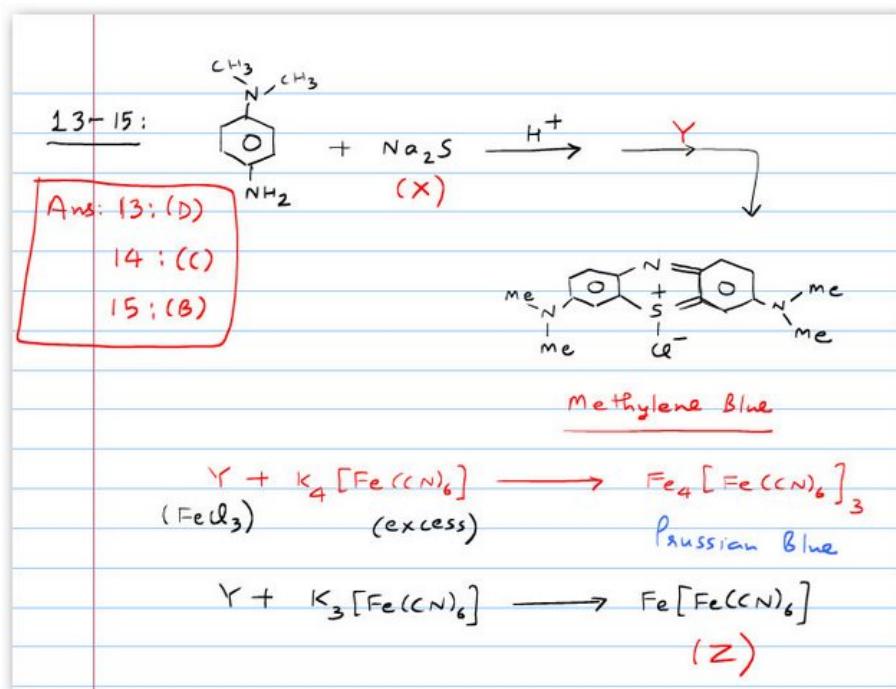
### Paragraph for Question Nos. 13 to 15

*p*-Amino-*N,N*-dimethylaniline is added to a strongly acidic solution of **X**. The resulting solution is treated with a few drops of aqueous solution of **Y** to yield blue coloration due to the formation of methylene blue. Treatment of the aqueous solution of **Y** with the reagent potassium hexacyanoferrate(II) leads to the formation of an intense blue precipitate. The precipitate dissolves on excess addition of the reagent. Similarly, treatment of the solution of **Y** with the solution of potassium hexacyanoferrate(III) leads to a brown coloration due to the formation of **Z**.

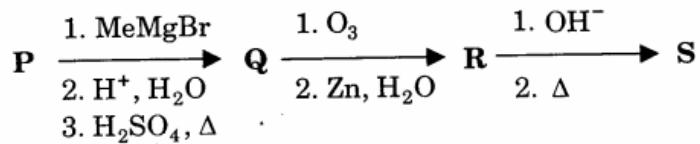
13. The compound **X** is

(A)  $\text{NaNO}_3$       (B)  $\text{NaCl}$       (C)  $\text{Na}_2\text{SO}_4$       (D)  $\text{Na}_2\text{S}$

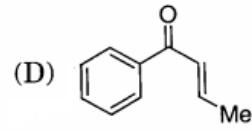
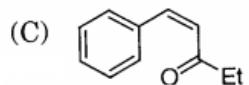
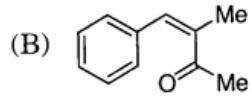
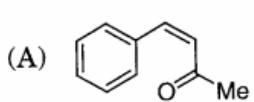
14. The compound **Y** is  
 (A)  $MgCl_2$       (B)  $FeCl_2$       (C)  $FeCl_3$       (D)  $ZnCl_2$
15. The compound **Z** is  
 (A)  $Mg_2[Fe(CN)_6]$       (B)  $Fe[Fe(CN)_6]$   
 (C)  $Fe_4[Fe(CN)_6]_3$       (D)  $K_2Zn_3[Fe(CN)_6]_2$



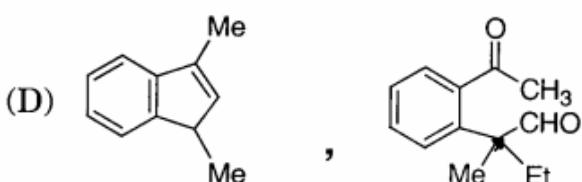
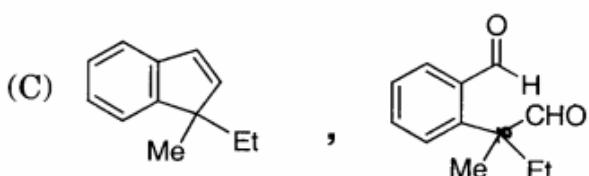
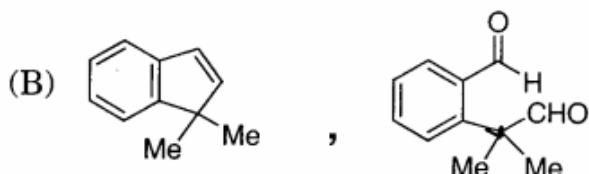
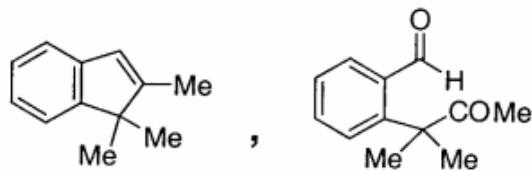
A carbonyl compound **P**, which gives positive iodoform test, undergoes reaction with  $\text{MeMgBr}$  followed by dehydration to give an olefin **Q**. Ozonolysis of **Q** leads to a dicarbonyl compound **R**, which undergoes intramolecular aldol reaction to give predominantly **S**.



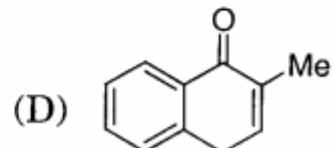
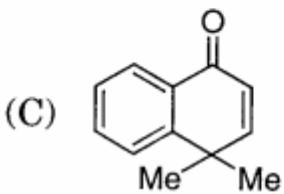
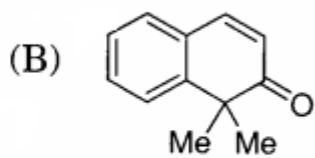
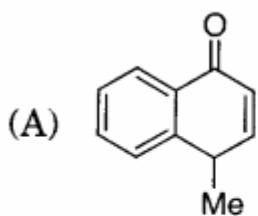
16. The structure of the carbonyl compound **P** is

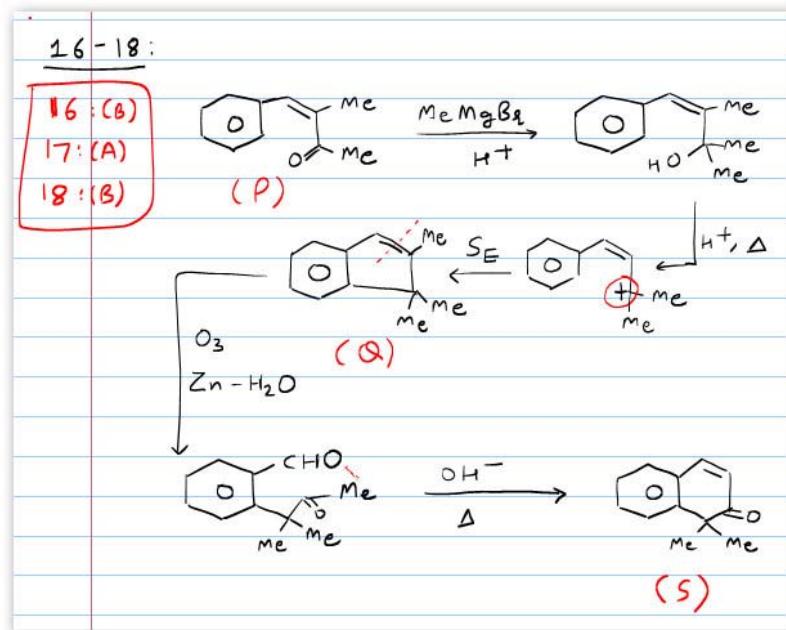


17. The structures of the products **Q** and **R**, respectively, are



18. The structure of the product **S** is





19. Match each of the diatomic molecules in **Column I** with its property/properties in **Column II**.

**Column I**

- (A)  $\text{B}_2$
- (B)  $\text{N}_2$
- (C)  $\text{O}_2^-$
- (D)  $\text{O}_2$

**Column II**

- (p) Paramagnetic
- (q) Undergoes oxidation
- (r) Undergoes reduction
- (s) Bond order  $\geq 2$
- (t) Mixing of 's' and 'p' orbitals

- 19: Ans: (A): p, q  
 (B): s  
 (C): p, q  
 (D): p, q, s

$B_2$  : B.O. = 1 (No. of unpaired e<sup>-</sup>s = 2)  
 ⇒ Paramagnetic

Adding one e<sup>-</sup> will go in bonding orbital.

$N_2$  : B.O. = 3

Adding/removing an e<sup>-</sup> will reduce its bond order.

In M.O. theory,

's' overlaps with 's'  
& 'p' overlaps with 'p'

$O_2^-$  : B.O. = 1.5  
 ↓ Paramagnetic  
 ↑

$O_2$  : B.O. = 2

will prefer the removal of e<sup>-</sup> from antibonding orbitals

20. Match each of the compounds in **Column I** with its characteristic reaction(s) in **Column II**.

**Column I**

- (A)  $CH_3CH_2CH_2CN$
- (B)  $CH_3CH_2OCOCH_3$
- (C)  $CH_3 - CH = CH - CH_2OH$
- (D)  $CH_3CH_2CH_2CH_2NH_2$

**Column II**

- (p) Reduction with Pd - C/H<sub>2</sub>
- (q) Reduction with SnCl<sub>2</sub>/HCl
- (r) Development of foul smell on treatment with chloroform and alcoholic KOH
- (s) Reduction with diisobutylaluminium hydride (DIBAL - H)
- (t) Alkaline hydrolysis

20:

(A) : q, s, t  
(B) : s, t  
(C) : p  
(D) : r

