

NAVODAYA VIDYALAYA SAMITI - HYDERABAD REGION

SECOND PRE BOARD EXAMINATION, 2022-23

CLASS -XII

SUBJECT- CHEMISTRY THEORY (043)

SET : 1


MARKING SCHEME

Section A (1m each)

1. A
2. C
3. C
4. B
5. A
6. D
7. A
8. D
9. D
10. B
11. D
12. B
13. C
14. B
15. D
16. C
17. B
18. C

Section B

19. O-Nitro phenol is steam volatile 1m
Reason : Intra molecular hydrogen bonding 1m

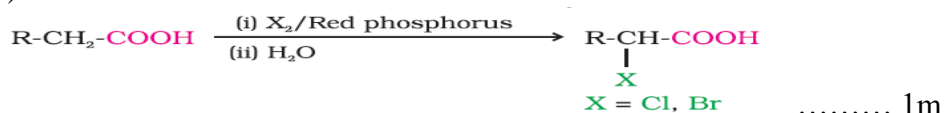
20. :  reacts faster by S_N1 mechanism 1m
it is a tertiary halide and it produces a stable tertiary carbocation...1m

21. $t_{99\%} = \frac{2.303}{k} \log (100/0.1)$ -----(1) 1/2m

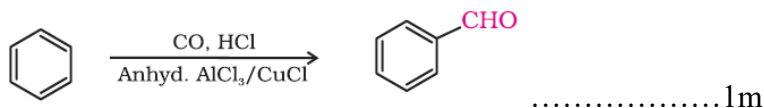
$$t_{1/2} = \frac{0.693}{k} \text{ -----(2) 1/2m}$$

$$t_{99\%} / t_{1/2} = 6.909 / 0.693 = 10 \text{ 1m}$$

22. A (i)

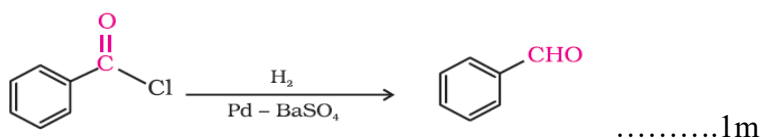


(ii)

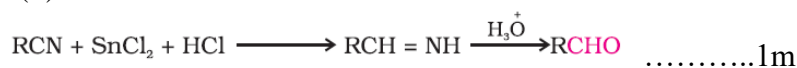


or

B. (i)



(ii)



23. According to Arrhenius equation,

$$\log k = \log A - \frac{E_a}{2.303 RT} \quad \dots\dots\dots 1/2\text{m}$$

$$\log k = 14.34 - 1.25 \times 10^4 \text{K}/T$$

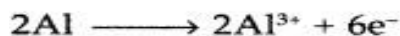
comparing both the equations:

$$\frac{E_a}{2.303 R} = 1.25 \times 10^4 \quad \dots\dots\dots 1/2\text{m}$$

$$E_a = 2.303 \times 8.314 \times 1.25 \times 10^4$$

$$= 239.34 \text{ kJ/mol} \quad \dots\dots\dots 1\text{m}$$

24



Hence, $n = 6$

$$\Delta G^\circ = -nFE_{\text{Cell}}^\circ$$

$$\Delta G^\circ = -6 \times 96500 \times 2.02 = -1169580 \text{ J mol}^{-1}$$

$$\therefore \Delta G^\circ = -116.958 \text{ KJ mol}^{-1}$$

$$\text{Now, } \Delta G^\circ = -2.303 RT \log K_c$$

$$\log K_c = \frac{-\Delta G^\circ}{2.303 RT} = \frac{1169580}{2.308 \times 8.314 \times 298}$$

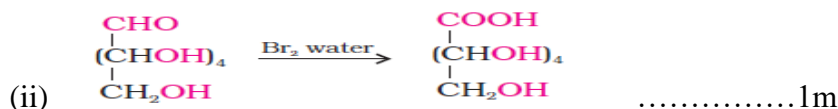
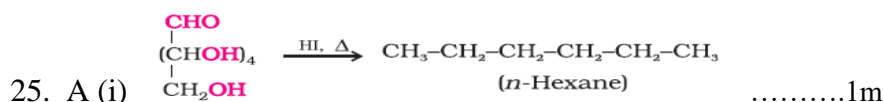
$$= \frac{1169580}{5705.84}$$

$$\therefore \log K_c = 205.009$$

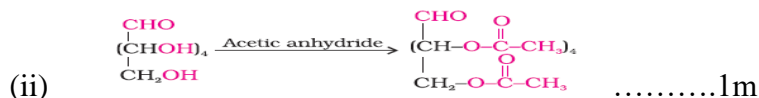
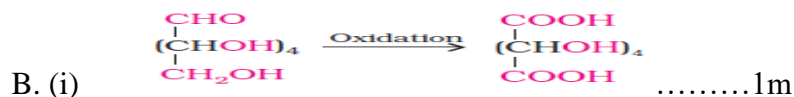
1 M

1 M

1 M

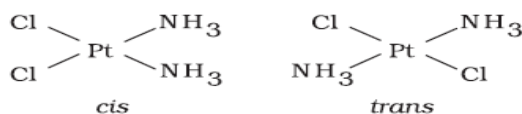


OR



Section C

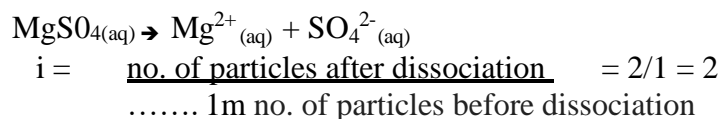
26. (i) structures of cis and trans isomers ($\frac{1}{2}$ m each)



(ii) Electronic configuration for d^4 ion if $\Delta_0 < P$ is $t_{2g}^3 e_g^1$ (high spin complex is formed).....1m

(iii) $[\text{Ni}(\text{H}_2\text{O})_6]^{2+}$ is an outer orbital complex due to weak field ligand H_2O and the presence of unpaired electrons undergoes d—d transition (by absorbing red light and shows green colour) while $[\text{Ni}(\text{CN})_4]^{2-}$ is an inner orbital complex and has no unpaired electrons hence colourless... 1m

27. Since MgSO_4 is an ionic compound, so undergoes complete ionisation in the following way:



$$\Delta T_b = i K_b m = i K_b \frac{W_2 \times 1000}{M_2 \times W_2} = 2 \times 0.52 \times \frac{4 \times 1000}{120 \times 100} = 0.346\text{K}$$

.....1m

Boiling point of water = 373.15 K

$$T_b = T_b^\circ + \Delta T_b = 373.15 \text{ K} + 0.346 = 373.496 \text{ K} \dots\dots\dots 1\text{m}$$

28. (a) Chlorobenzene has lower dipole moment than cyclohexyl chloride due to lower magnitude of -ve charge on the Cl atom and shorter C – Cl distance. Due to greater S-character, a sp^2 -hybrid carbon is more electronegative than a sp^3 -hybrid carbon. Therefore, the sp^2 -hybrid carbon of C – Cl bond in chlorobenzene has less tendency to release electrons to Cl than a sp^3 hybrid carbon of cyclohexyl chloride. 1m

(b) Alkyl halides and polar molecules are held together by dipole-dipole interaction. The molecules of H_2O are held together by H- bonds. Since the new forces of attraction between water and alkyl halide molecules are weaker than the forces of attraction already existing between alkyl halide-alkyl halide molecules and water- water molecules, therefore alkyl halides are immiscible (not soluble) with water. 1m

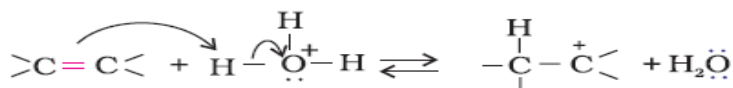
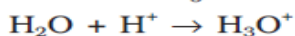
(c) n-Butyl bromide has higher boiling point than 3^o-Butyl bromide because it has larger surface area hence have more VanderWaals' forces. 1m

29. a. CH₃—NH₂ has higher boiling point than (CH₃)₃N. (1/2m)
Reason: CH₃—NH₂ shows strong intermolecular H bonding whereas (CH₃)₃N has weak dipole dipole forces. 1/2m

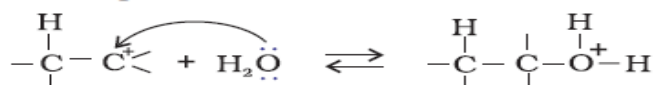
- b. A = CH₃-CH₂-CN,1/2 m
Propanenitrile1/2m
B = CH₃-CH₂-CH₂-NH₂,1/2 m
Propanamine1/2 m

30.

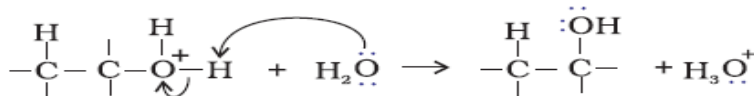
Step 1: Protonation of alkene to form carbocation by electrophilic attack of H₃O⁺.



Step 2: Nucleophilic attack of water on carbocation.



Step 3: Deprotonation to form an alcohol.



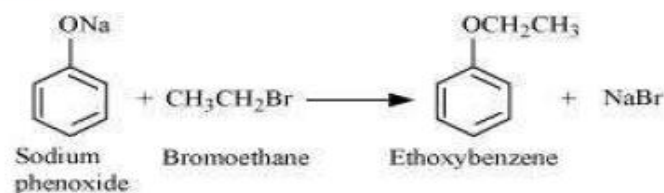
mechanism with three steps: 1m for each step

OR

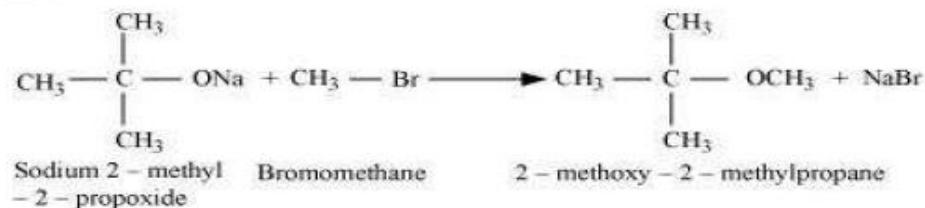


(i) Sodium propoxide 1-Bromopropane 1- Propoxypropane

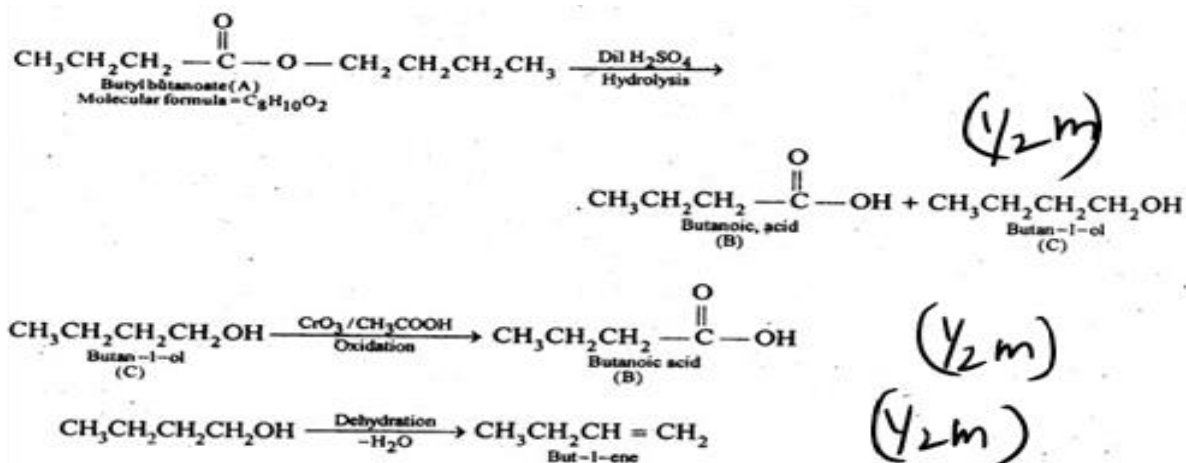
(ii)



(iii)



Three equations : 1m for each



b. (i) Iodoform test :Propanal has an aldehydic functional group and propanone is a methyl ketone. Propanal on reaction with sodium hypoiodite does not form a yellow coloured precipitate. But propanone on reaction with sodium hypoiodite form a yellow coloured precipitate 1/2 m

Relevant Chemical equation1/2 m

Or Tollen's reagent, Benedict solution and Fehling solution are all used to distinguish between aldehydes and ketones. Propanal($\text{CH}_3\text{CH}_2\text{CHO}$) is an aldehyde and thus gives positive test with all the three reagents whereas propanone(CH_3COCH_3) is a ketone and thus does not give any results with the three reagents.....1/2 m

Relevant chemical equation1/2 m

(ii) Phenol gives violet colour with neutral FeCl_3 while benzoic acid gives buff colored ppt. Benzoic acid react with NaHCO_3 and form CO_2 gas ,But does not react with NaHCO_31/2 m

Relevant chemical equation1/2 m

34 a. $m = 1.5\text{g}$, $I = 1.5\text{A}$, Molar mass = 108 g/mol ,
 $1F = 96500\text{C/mol}$
 $m = zIt$ 1/2 m
 m
 $1.5 = \frac{108}{1 \times 96500} \times 1.5 t$ 1m
 $t = 893\text{s}$ 1/2 m

b. $\text{Fe} \longrightarrow$
 $\text{Fe}^{2+} + 2e$
 $\text{Fe} + 2\text{H}^+ \longrightarrow \text{Fe}^{2+} + \text{H}_2$
 $n = 2$ 1/2m
 $E^0_{\text{cell}} = 0.00 - (-0.44) = 0.44\text{V}$ 1/2m
 $E_{\text{cell}} = E^0_{\text{cell}} - \frac{0.0591}{2} \log[\text{Fe}^{2+}]$ 1/2m
 $= 0.44 - \frac{0.0591}{2} \log[10^{-3}]$
 $= 0.4105\text{V}$ 1/2m

c. Primary cell. It has more efficiency and its voltage remains constant over a longer period of time.1m

35. (a) In transition elements, there are large number of unpaired electrons in their atoms, thus they have a stronger inter atomic interaction and thereby stronger bonding between the atoms. Due to this they have high enthalpies of atomization..... 1m

(b) Due to variable valency and the availability of d-orbitals, they can easily form intermediate products which are converted to products. The sizes of transition metal atoms and ions are also favourable for transition complex formation with the reactants. They can adsorb reactants on their surface and catalyze reactions. 1m

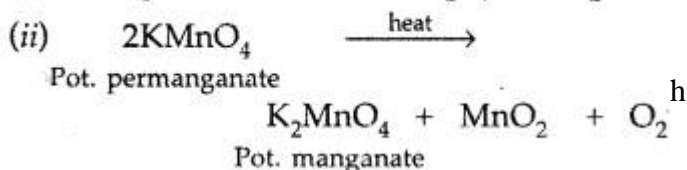
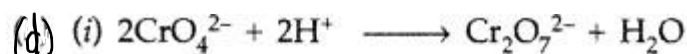
(c) Divalent ion with atomic number

26 is Fe^{2+} Electronic conf of Fe^{2+} -

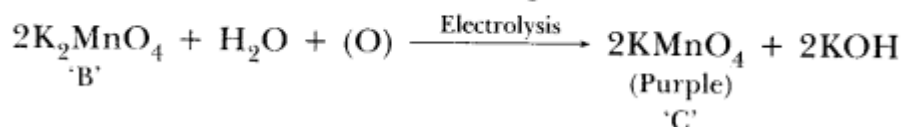
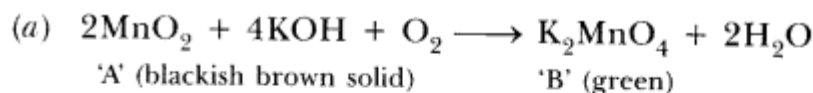
$[\text{Ar}] 3d^6$

No. of unpaired electron (n) = 4

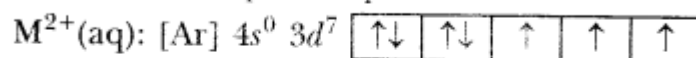
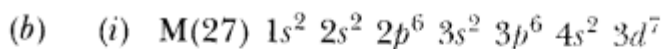
$$\mu = \sqrt{n(n+2)} = \sqrt{24} = 4.9 \text{ BM} \dots\dots\dots 1\text{m}$$



(II)



..... 3m



$$\mu = \sqrt{n(n+2)}$$

$$\mu = \sqrt{3 \times 5} = \sqrt{15} = 3.87 \text{ BM.}$$

.....1m

(ii) Cr is typically hard metal due to presence of large number of unpaired electrons, metal-metal interaction is strong whereas mercury does not have unpaired electrons and has large size, therefore, forms weak metallic bond. 1m
