SAMPLE PAPER SOLUTION (2021-22)

CLASS XII TERM – II

CHEMISTRY THEORY (043)

Important Instructions

- 1. This is only a suggestive answer key/marking scheme.
- 2. Any other correct response/(s) are also accepted.

| Q.No. | Suggestive Answers | Step Marking |
|-------------------|---|--------------|
| 1. a. | This is due to stearic hindrance/Nu has difficulty in approaching carbon | 1 |
| b. | The NH ₂ in conjugation with carbonyl involves in RESONANCE, loses its nucleophilicity | 1 |
| 2. a. | Increases linearly | 1/2 |
| b. | Increases steeply, the degree of dissociation is higher in weak electrolyte | 1/2 + 1 |
| 3. a | The negative charge is dispersed more on, more electronegative atom i.e. Oxygen in carboxylate ion. | 1 |
| 3. b. | Iodoform test/Fehling's test/Tollens' test | 1 |
| 4. a | Because of lower electronegativity of Nitrogen than oxygen/The removal of H as H^{\dagger} is weaker. | 1 |
| b. | Primary amines forms more number of Intermolecular H bond/Stronger intermolecular forces of attraction | 1 |
| С. | In Aromatic amines lone pairs of nitrogen does resonance with benzene, electron density decreases. | 1 |
| OR 4.a. | Due to resonance, aryl part does not detach with halide and can't be introduced on nitrogen salt of Phthalimide. | 1 |
| b. | In aromatic salt, Due to resonance, diazonium group does not leave easily and gets extra stability | 1 |
| с. | Nitrogen has limited covalency/due to absence of d orbitals | 1 |
| 5.a. | $[Co(NH_3)_4(H_2O)Cl]Cl_2$ | 1 |
| b. | [NiCl ₄] ^{2–} contains unpaired electrons | 1 |
| с. | $[Co(NH_3)_6]^{3+}$, Co is $d^2 sp^3$ and $[Ni(NH_3)_6]^{2+}$, Ni is $sp^3 d^2 hybridised$. | 1 |
| OR a. | If $\Delta o > P$, Pairing occurs and if $\Delta o < P$, pairing does not occur in d subshell. | 1 |
| b. | i) $t_2^4 e_g^2$ ii) $t_2^6 e_g^0$ | 1+1 |
| 6.a. | i) $t_2^4 e_g^2$ ii) $t_2^6 e_g^0$ According to CFT Cr^{3+} is more stable in d3 (half filled) system than in d^5 , there Cr loses one electron to be oxidized and behave as reducing agent. | 1 |
| b. | This is due to higher ionization enthalpy of Cu as it is a contributory factor to Ev. | 1 |
| с. | Mn in +2 state is d5 and this state is highly stable. Conversion from d ⁵ to d ⁴ is highly unstable. | 1 |

| 7. | $2H_2O + 4NH_3 + 2Ag + $ Silver mirror C_2H_5 NO_2 COO^{-} C_2H_5 | 1+1+1 (one mark for each correct reaction) |
|------------|---|--|
| | $CH = NNH$ NO_2 $2, 4-Dinitrophenyl hydrazine$ $+ H_2O \leftarrow$ $2, 4-DNP$ derivative | |
| 8.a. | On increasing temperature molecules on the surface of adsorbent aquires energy ans desorb. | 1 |
| b. | This is due to higher surface area. | 1 |
| с. | Because energy releases when bond of adsorbate is formed with adsorbent. | 1 |
| 9 | (<i>i</i>) $CH_{3}COOH \xrightarrow{SOCl_{2}}{-SO_{2}, -HCl} CH_{3}COCI \xrightarrow{NH_{3}(excess)}{NH_{4}Cl} CH_{3}CONH_{2} \xrightarrow{Br_{7}/NaOH} CH_{3}NH_{2}$ (<i>ii</i>) $CH_{3}(CH_{2})_{4}CN \xrightarrow{H_{3}O^{+}}{CH_{3}(CH_{2})_{4}COOH} \xrightarrow{SOCl_{2}}{-SO_{2}, -HCl} CH_{3}(CH_{2})_{4}COCI \xrightarrow{NH_{3}(excess)}{-NH_{4}Cl} CH_{3}(CH_{2})_{4}CONH_{2}$ $CH_{3}(CH_{2})_{4}NH_{2} \xleftarrow{Br_{2}/NaOH} CH_{3}(CH_{2})_{4}NH_{2}$ (<i>iii</i>) $CH_{3}OH \xrightarrow{PCl_{5}}{-POCl_{3}} CH_{3}CI \xrightarrow{KCN(alc)}{CH_{3}CN} CH_{3}CN \xrightarrow{H_{3}O^{+}}{CH_{3}COOH}$ | 1+1+1 (one mark for each correct reaction) Or any other correct conversion |
| OR 9.a. | $NH_3 (alc) + C_2H_5CI \longrightarrow C_2H_5NH_2 + HCI$ | 1 |
| b. | $C_6H_5NC + H_2O + KCI$ | 1 |
| с. | $p-NH_3^+C_6H_4SO_3^-$ | 1 |
| 10. | $E_{cell} = E_{cell}^{o} - \frac{0.0591}{n} \log \frac{[Ni^{2+}]}{[Ag^{+}]^{2}}$ | 1/2 |
| | $= 1.05V - \frac{0.0591}{2} \log \frac{0.160}{(0.002)^2}$ | 1/2 |
| | $= 1.05 - \frac{0.0591}{2} \log(4 \times 10^4)$ = 1.05 - $\frac{0.0591}{2} (4.6021)$ = 1.05 - 0.14 V | 1 |
| | = 1.05 - 0.14 V = 0.91 V | $\frac{1}{2} + \frac{1}{2}$ |
| 11.a. | This is due to high electronegativity and small size | 1 |
| b. | Ce ⁴⁺ . Tr ⁴⁺ , Pr ⁴⁺ (Any Two) | 1 |
| с. | Shielding of 5f is more poorer than 4f. | 1 |
| OR a. | This is due to electron – electron repulsion at the later stage of the series. | 1 |
| b. | Lanthanoid Contraction. | 1 |
| с. | It is due to lowest intermolecular forces of attraction in zinc as it is d^{10}/no unpaired electrons. | 1 |
| 12.a. | (i) 0.2M (ii) 0.1M | 1+1 |

| b. | Rate = $[A]^{\alpha}$ $[B]^{\beta}$ $r_1 = 5.07 \times 10^{-5} = (0.20)^{\alpha} (0.30)^{\beta}$ (i) | 1/2 1/2 |
|----|---|------------|
| | $r_2 = 5.07 \times 10^{-5} = (0.20)^{\alpha} (0.10)^{\beta} \qquad(ii)$ $r_3 = 1.43 \times 10^{-4} = (0.40)^{\alpha} (0.05)^{\beta} \qquad(iii)$ | /2 1⁄2 |
| | $\frac{r_1}{r_2} = \frac{5.07 \times 10^{-5}}{5.07 \times 10^{-5}} = \frac{(0.2)^{\alpha}}{(0.2)^{\alpha}} \frac{(0.30)^{\beta}}{(0.10)^{\beta}} = 1 = (3)^{\beta}$ $\therefore \beta = 0$ | 1/2 |
| | $\frac{r_3}{r_2} = \frac{1.43 \times 10^{-4}}{5.06 \times 10^{-5}} = \frac{(0.40)^{\alpha}}{(0.20)^{\alpha}} \frac{(0.05)^{\beta}}{(0.10)^{\beta}}$ | 1/2 |
| | $= 2.826 = 2^{\alpha} \left(\frac{1}{2}\right)^{\beta}$ 2.826 = 2 ^{\alpha} (as \beta = 0) Taking log on both the sides, we get log 2.826 = \alpha log 2 \approx 0.4511 = \alpha \times 0.3010 | |
| | $\alpha = \frac{0.4511}{0.3010} = 1.498 = 1.5$ | 1/2 |
| | \therefore Order with respect to A = 1.5 and order with respect to B = 0 | |
| | | |