

DAV PUBLIC SCHOOLS, ODISHA ZONE

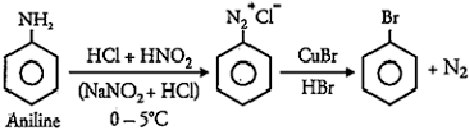
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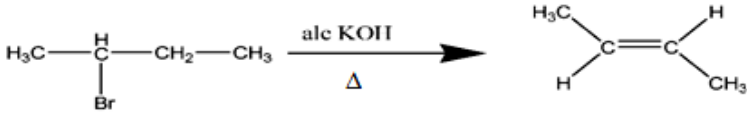
CLASS : XII , SUBJECT : CHEMISTRY

MARKING SCHEME(SET-2)

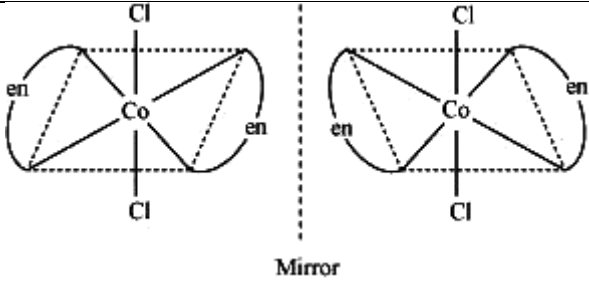
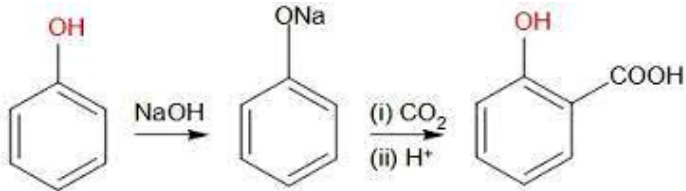
QSN O	Value Points	Marks Allotted	PAGE NO. OF NCERT /TEXT BOOK
1	a) He	1	45
2	a)The activation energy of forward reaction is $E_1 + E_2$ and the product is less stable than reactant.	1	113
3	a) Sc	1	311
4	a) But-3-en-2-ol	1	289
5	a) $0.005 \text{ molL}^{-1}\text{s}^{-1}$	1	96
6	c) Nearly same atomic size	1	213
7	c) 16 times	1	100
8	c) $8,000 \text{ cm}^{-1}$	1	252
9	c) 3	1	226
10	a)2-Methylpropene	1	206
11	a) $\text{Cr} > \text{Mn} > \text{V} > \text{Ti}$	1	221
12	d) benzyl alcohol	1	222
13	b) Both A and R are true but R is not the correct explanation of A.	1	46
14	d) A is false but R is true.	1	210
15	b) Both A and R are true but R is not the correct explanation of A.	1	101
16	b) Both A and R are true but R is not the correct explanation of A.	1	295
17	<p>The net reaction is :</p> $\text{Ni}(s) + 2\text{Ag}^+(aq) \rightarrow \text{Ni}^{2+}(aq) + 2\text{Ag}(s)$ <p>According to Nernst equation,</p> $E_{\text{cell}} = E_{\text{cell}}^{\circ} - \frac{0.0591}{n} \log \frac{[\text{Anode}]}{[\text{Cathode}]}$ $E_{\text{cell}} = \left[E_{\text{Ag}^+ / \text{Ag}}^{\circ} - E_{\text{Ni}^{2+} / \text{Ni}}^{\circ} \right] - \frac{0.0591}{2} \log \frac{[\text{Ni}^{2+}]}{[\text{Ag}^+]^2}$ $= 1.05 - \frac{0.0591}{2} \log \frac{0.16}{(0.002)^2}$ $= 1.05 - \frac{0.0591}{2} \log(4 \times 10^4) = 1.05$ $- \frac{0.0591}{2} (\log 4 + 4 \log 10)$ $= 1.05 - \frac{0.0591}{2} \times 4.6020 = 1.05 - 0.1360$ $= 0.914V$	<p>$\frac{1}{2}$</p> <p>$\frac{1}{2}$</p> <p>$\frac{1}{2}$</p> <p>$\frac{1}{2}$</p>	70

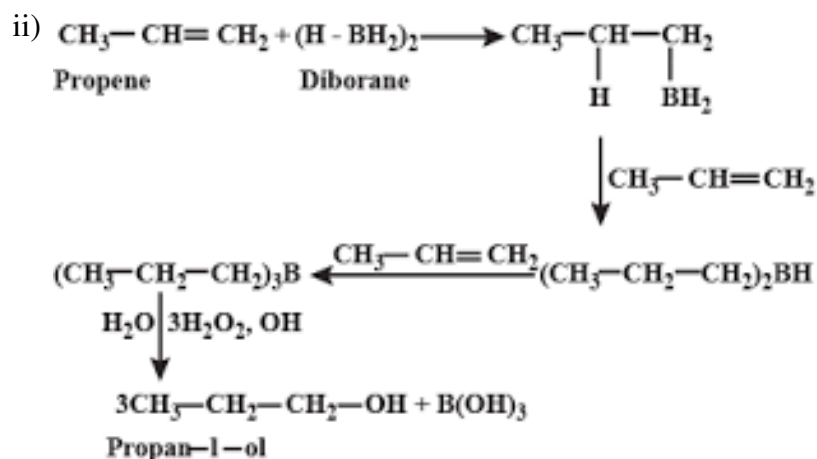
18	a) $6 \text{Fe}^{2+} + \text{Cr}_2\text{O}_7^{2-} + 14 \text{H}^+ \rightarrow 2 \text{Cr}^{3+} + 6 \text{Fe}^{3+} + 7 \text{H}_2\text{O}$ b) $2\text{MnO}_4^- + 5\text{C}_2\text{O}_4^{2-} + 16\text{H}^+ \rightarrow 2\text{Mn}^{2+} + 10\text{CO}_2 + 8\text{H}_2\text{O}$	1 1	226																
19	A= C_2H_4 , B= CH_3CH_3 , C= $\text{CH}_3\text{OC}_2\text{H}_5$, D= $\text{CH}_3\text{CH}_2\text{NC}$	$\frac{1}{2} \times 4$	299-310																
20	a) <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th style="width: 30%;">compounds</th> <th style="width: 30%;">Tests/reagents</th> <th style="width: 40%;">observation</th> </tr> </thead> <tbody> <tr> <td>phenol</td> <td rowspan="2">Neutral FeCl_3 solution</td> <td>Violet colour solution</td> </tr> <tr> <td>ethanol</td> <td>No such obs.</td> </tr> </tbody> </table> b) <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th style="width: 30%;">compounds</th> <th style="width: 30%;">Tests/reagents</th> <th style="width: 40%;">observation</th> </tr> </thead> <tbody> <tr> <td>tert-butyl alcohol</td> <td rowspan="2">Lucas test (conc.HCl+anh.ZnCl₂)</td> <td>Turbidity occurs immediately.</td> </tr> <tr> <td>n- butyl alcohol</td> <td>No such obs.</td> </tr> </tbody> </table>	compounds	Tests/reagents	observation	phenol	Neutral FeCl_3 solution	Violet colour solution	ethanol	No such obs.	compounds	Tests/reagents	observation	tert-butyl alcohol	Lucas test (conc.HCl+anh.ZnCl ₂)	Turbidity occurs immediately.	n- butyl alcohol	No such obs.	$\frac{1}{2} + \frac{1}{2}$ $\frac{1}{2} + \frac{1}{2}$	341 338
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21	a) The enthalpies of atomization of a transition metal are high because they have a large number of unpaired electrons resulting strong inter atomic metallic bonding. b) Weak inter atomic metallic bonding due to absence of unpaired electron. OR a) In Comparison to Fe^{2+} , Cr^{2+} is a stronger reducing agent because in formation of Cr^{3+} from Cr^{2+} changes is from $d^4 \rightarrow d^3$. In d^3 electronic configuration t_{2g} orbitals are half filled. But in Fe^{2+} to Fe^{3+} Changes is d^6 to d^5 b) Atomic number (Z)=27, it is Co with configuration $3d^7, 4s^2$ In Co^{2+} , the configuration is $3d^7$. Now, Number of unpaired electrons =3 magnetic moment, $\mu = \sqrt{n(n+2)} = \sqrt{3(3+2)} = 3.87 \text{ BM}$	1 1 1 1	218																
22	a) It states that "the partial pressure of the gas in vapour phase (p) is proportional to the mole fraction of the gas (x) in the solution". b) Since number of particles decreases, hence van't Hoff factor (i) will decrease and freezing point of the solution will increase. c) Molality is considered better for expressing the concentration as compared to molarity because the molarity changes as volume of the solution changes with temperature.	1 1 1	46 58 37																
23	a) $\text{Al}_2\text{O}_3 + 6e^- \longrightarrow 2\text{Al} + 3\text{O}^{2-}$ $\text{6F} \qquad (2 \times 27) \text{ g}$ <p>To produce 54 g of Al, charge needed = 6F</p> <p>To produce 40 g of Al, charge needed = $\frac{(40 \text{ g})}{(54 \text{ g})} \times (6F) = 4.44F$.</p> b) At anode= Br_2 , at cathode= Cu c) Δ° for NaBr is calculated by the following expression. $\Delta^\circ \text{NaBr} = \lambda^\circ \text{NaCl} + \lambda^\circ \text{KBr} - \lambda^\circ \text{KCl}$ $= 126 + 152 - 150 = 128 \text{ Scm}^2 \text{ mol}^{-1}$	1 $\frac{1}{2} + \frac{1}{2}$ 1	94 87 83																

24	<p>a) Bis(ethane 1,2-diamine) dihydroxidochromium(III) chloride</p> <p>b) In $[\text{Ni}(\text{H}_2\text{O})_6]^{2+}$, water is a weak field ligand. Therefore, there are unpaired electrons in Ni^{2+}. In this complex, the d electrons from the lower energy level can be excited to the higher energy level i.e., the possibility of d-d transition is present. Hence, $[\text{Ni}(\text{H}_2\text{O})_6]^{2+}$ is coloured.</p> <p>In $[\text{Ni}(\text{CN})_4]^{2-}$, the electrons are all paired as CN^- is a strong field ligand. Therefore, d-d transition is not possible in $[\text{Ni}(\text{CN})_4]^{2-}$. Hence, it is colourless.</p> <p>c) $t_{2g}^4 e_g^0$</p>	1 1 1	249 259
25	<p>a)</p> $\log k = \log A - \frac{E_a}{2.303RT}$ $14.2 - \frac{1.0 \times 10^4}{T} = \log A - \frac{E_a}{2.303RT}$ $1.0 \times 10^4 = \frac{E_a}{19.14}$ $E_a = 19.14 \times 10^4 \text{ J mol}^{-1} = 191.4 \text{ KJ mol}^{-1}$ <p>For first order</p> $k = 0.693/t_{1/2}$ $= 0.693/200 \text{ min}^{-1}$ $= 0.3465 \text{ min}^{-1}$ <p>b) elementary reaction.</p>	1/2 1/2 1/2 1	113
26	<p>Any three</p> <p>a)</p>  <p>b) As all the hydrogen atoms are equivalent and replacement of any hydrogen will give the same product.</p> $ \begin{array}{c} \text{CH}_3 \\ \\ \text{H}_3\text{C} - \text{C} - \text{CH}_3 \\ \\ \text{CH}_3 \end{array} \quad (\text{Neopentane}) $ <p>c) In haloarenes, NO_2 group present at o/p position results in the stabilisation of resulting carbanion by -R and -I effects and therefore increases the reactivity of haloarenes towards nucleophilic substitution reactions.</p> <p>d) 2, 2-Bis (4-chlorophenyl)-1,1,1-trichloroethane</p>	1 1 1 1	206 301 313 318
27	<p>a) $\text{CH}_3\text{COOC}_2\text{H}_5 + \text{H}_2\text{O} \xrightarrow{\text{H}^+} \text{CH}_3\text{COOH} + \text{C}_2\text{H}_5\text{OH}$ (any other suitable example)</p>	1	102

	b) $-\frac{dx}{dt} = k[A][B]^2$	1	
	c) 1	1	
28	$A = \pi r^2 = 3.14 \times (0.5)^2 \text{ cm}^2 = 0.785 \text{ cm}^2$ $\rho \text{ (resistivity)} = \frac{R \times A}{l} = \frac{5.55 \times 10^3 \Omega \times 0.785 \text{ cm}^2}{50 \text{ cm}} = 87.135 \Omega \text{ cm.}$ $\kappa = \frac{1}{\rho} = \frac{1}{87.135 \Omega \text{ cm}} = 0.01148 \text{ S cm}^{-1}$ $\Lambda_m = \frac{\kappa \times 1000}{M} = \frac{0.01148 \times 1000}{0.05 \text{ M}} = 229.6 \text{ S cm}^2 \text{ mol}^{-1}$	1 1 1	242
29	a) inversion of configuration b)  c) But-2-ene  OR a) 1-Bromopentane will be more reactive as it is least crowded for an SN2 reaction. b) 2-Bromopentane has a chiral carbon. Therefore, it is optically active.	1 1 1+1 1 1	171-174
30	a) $\text{Zn(s)} \mid \text{Zn}^{2+}(\text{aq}) \parallel \text{Cu}^{2+}(\text{aq}) \mid \text{Cu (s)}$ b) Once the salt bridge is removed from the cell, the voltage drops to zero. No flow of current. OR The cell will behave as an electrolytic cell. c) $Q = 1.5 \times 900C = 1350C$. 96500 C deposits 1 g eq. hence, $E = \frac{0.783}{1350} \times 96500 = 55.97$	1 1 1 1	66,86
31	a) i) The bonds between chloroform molecules and molecules of acetone are dipole-dipole interactions but on mixing, the chloroform and acetone molecules start forming hydrogen bonds which are stronger bonds resulting in the release of energy. This gives rise to an increase in temperature. ii) To avoid bends, as well as, the toxic effects of high concentrations of nitrogen in the blood, the tanks used by scuba divers are filled with air diluted with helium. iii) The magnitude of osmotic pressure is large even for very dilute solution and it can be measured at room temperature. (any other suitable reason)	1 1 1	43,45,54

	<p>b)</p> $M_B = \frac{K_f \times W_B \times 1000}{\Delta T_f \times W_A}$ $M_B = \frac{3.83 \text{ K kg mol}^{-1} \times 2.56 \times 1000 \text{ g kg}^{-1}}{100 \text{ g} \times 0.383 \text{ K}} = 256 \text{ g mol}^{-1}$ <p>Now, molecular mass of $S_x = x \times 32 = 256$</p> $x = \frac{256}{32} = 8$ <p>Therefore, formula of sulphur = S_8</p>	<p>$\frac{1}{2}$</p> <p>$\frac{1}{2}$</p> <p>$\frac{1}{2}$</p> <p>$\frac{1}{2}$</p>	
	<p style="text-align: center;">OR</p> <p>a) i) Beaker 1: Hypotonic solution, Beaker 2: Hypertonic solution ii) In beaker 3 the size of potato cube remains the same because of isotonic solution which has the same concentration of solutes as that of potato cells. So water is neither lost or gained by the potato cells.</p> <p>b) $\Delta T_f = 0 - (-0.068) = 0.068 \text{ K}$ $\Delta T_f = i \times K_f \times m$ $0.068 = i \times 1.86 \times 0.01$ So, $i = 3.6559$ Again, $\alpha = \frac{i-1}{n-1}$ for $\text{AlCl}_3 \rightleftharpoons \text{Al}^{3+} + 3\text{Cl}^-$; $n = 1 + 3 = 4$ $\alpha = \frac{3.6559 - 1}{4 - 1} = 0.8833$ % of dissociation = 88.33%</p> <p>c) The freezing point of water decreases, due to which the snow on the road starts to melt and clears the road.</p>	<p>$\frac{1}{2} + \frac{1}{2}$</p> <p>1</p> <p>$\frac{1}{2}$</p> <p>$\frac{1}{2}$</p> <p>$\frac{1}{2}$</p> <p>1</p>	51,54
32	<p>Any five:</p> <p>a) Hybridization: d^2sp^3, Magnetic character: Diamagnetic</p> <p>b) $[\text{Cr}(\text{H}_2\text{O})_5\text{Cl}]\text{Cl}_2 \cdot \text{H}_2\text{O}$</p> <p>c) No, ionization isomers are possible by exchange of ligand with counter ion only and not by exchange of central metal ion.</p> <p>d) In both $[\text{NiCl}_4]^{2-}$ and $[\text{Ni}(\text{CN})_4]^{2-}$, Ni is in +2 oxidation state with configuration $3d^8$ and it contains two unpaired electrons. In $[\text{NiCl}_4]^{2-}$ due to presence of weak ligand Cl^- no pairing takes place and hence it is paramagnetic whereas in $[\text{Ni}(\text{CN})_4]^{2-}$, CN^- is a strong field ligand and pairing occurs and hence it becomes diamagnetic.</p> <p>e) $[\text{Co}(\text{NH}_3)_6]\text{Cl}_3 > [\text{Cr}(\text{NH}_3)_5\text{Cl}]\text{Cl}_2 > [\text{Co}(\text{NH}_3)_4\text{Cl}_2]\text{Cl} > [\text{Co}(\text{NH}_3)_3\text{Cl}_3]$</p> <p>f) Trans isomer is optically inactive as it has a plane of symmetry</p>	<p>$\frac{1}{2} + \frac{1}{2}$</p> <p>1</p> <p>$\frac{1}{2} + \frac{1}{2}$</p> <p>1</p> <p>1</p> <p>$\frac{1}{2} + \frac{1}{2}$</p>	244, 249, 252

	 <p style="text-align: center;">Mirror</p> <p>g) The central atom is an electron pair acceptor so it is a Lewis acid.</p>	1	
33	<p>a)</p> $\begin{array}{ccc} \text{CH}_3\text{---CH---CH}_2\text{OH} & \xrightarrow[\text{H}_2\text{SO}_4]{\text{K}_2\text{Cr}_2\text{O}_7} & \text{CH}_3\text{---CH---CO}_2\text{H} \\ & & \\ \text{CH}_3 & & \text{CH}_3 \\ \text{(A) 2-Methylpropanol} & & \text{(B) 2-Methylpropanoic acid} \end{array}$ $\begin{array}{ccc} \downarrow \text{H}_2\text{SO}_4/\text{Heat} & & \\ \text{CH}_3\text{---C=CH}_2 & \xrightarrow[\text{H}_2\text{O}/\text{H}^+]{\text{H}_2\text{SO}_4} & \begin{array}{c} \text{OH} \\ \\ \text{CH}_3\text{---C---CH}_3 \\ \\ \text{CH}_3 \end{array} \\ \text{(C) 2-Methylpropene} & & \text{(D) 2-Methyl-2-propanol} \end{array}$ <p>A = $\begin{array}{c} \text{CH}_3\text{---CH---CH}_2\text{OH} \\ \\ \text{CH}_3 \end{array}$ B = $\begin{array}{c} \text{CH}_3\text{---CH---CO}_2\text{H} \\ \\ \text{CH}_3 \end{array}$</p> <p>C = $\begin{array}{c} \text{CH}_3\text{---C=CH}_2 \\ \\ \text{CH}_3 \end{array}$ D = $\begin{array}{c} \text{OH} \\ \\ \text{CH}_3\text{---C---CH}_3 \\ \\ \text{CH}_3 \end{array}$</p> <p>b) The commercial alcohol is made unfit for drinking by mixing in it some copper sulphate (to give it a colour) and pyridine (a foul smelling liquid). It is known as denaturation of alcohol.</p>	<p>$\frac{1}{2}$</p> <p>$\frac{1}{2}$</p> <p>1</p> <p>$\frac{1}{2} \times 4 = 2$</p> <p>1</p>	<p>210-216</p>
OR			
	<p>a)</p> <p>i)</p> 	1	200, 213, 215

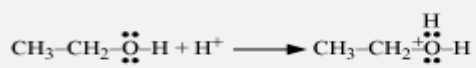


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b) Ethoxyethane

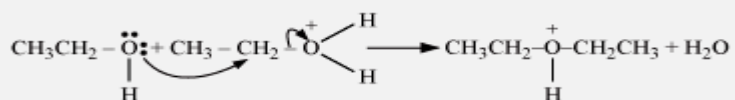
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Step 1



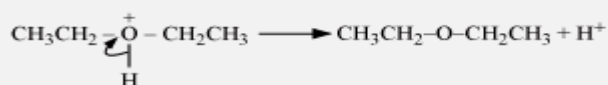
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Step 2



1

Step 3



1
