MARKING SCHEME

SAMPLE PAPER 1

SECTION A

Q.No.	Value Point	Marks
1(i)	D	1
(ii)	В	
	OR	1
	Α	
(iii)	Α	1
(iv)	С	1
2(i)	В	1
(ii)	Α	1
(iii)	Α	1
(iv)	Α	
	OR	1
	В	
3	С	1
4	D	
	OR	1
	С	
5	С	1
6	В	
	OR	1
	В	
7	В	
	OR	1
	D	
8	А	
	OR	1
	Α	
9	С	1
10	Α	1
11	Α	1
12	Α	1
13	D	1
14	В	
	OR	1
	В	
15	В	1
16	Α	1

SECTION B, C, D

Q.No.	VALUE POINTS	MARKS	
Q.110.	SECTION B		
17	Nitro group at ortho position withdraws the electron density from the benzene ring and thus facilitates the attack of the nucleophile on haloarene.	2	
	$\begin{array}{c} Cl & O \\ OH + \end{array} \xrightarrow{\begin{tabular}{c} Cl & OH \\ \hline & & & \\ \hline \hline & & & \\ \hline & & & \\ \hline \hline & & & \\ \hline \hline \\ \hline & & & \\ \hline \hline \\ \hline & & & \\ \hline \hline \\ \hline \hline & & & $		
	$= \underbrace{\overset{\operatorname{Cl}}{\bigoplus} \overset{\operatorname{OH}}{\bigoplus} \overset{\operatorname{OH}}{\bigoplus} }_{\operatorname{Fast step}} \xrightarrow{\operatorname{OH}} \underbrace{\overset{\operatorname{OH}}{\bigoplus} \overset{\operatorname{OH}}{\bigoplus} }_{\oplus \operatorname{O}} + \operatorname{Cl}^{\ominus}$		
	OR		
	(i) NH_2 N_2Cl Cl Cl N_2Cl Cl Cl N_2Cl Cl Cu_2Cl_2	1	
	273 – 278 K		
	(ii) $CH_3CH(Br)CH_3 \xrightarrow{\text{alc KOH}} CH_3CH=CH_2 \xrightarrow{\text{HBr, organic peroxide}} HBr, organic peroxide$	1	
18	$\Delta Tb = K_b m \Delta Tb = 101.04-100 = 1.04 \text{ °C}$ or m= 1.04 /0.52 = 2 m 2 m solution means 2 moles of solute in 1 kg of solvent.	1	
	2 m aq solution of urea means 2 moles of urea in 1kg of water. No. of moles of water = $1000/18 = 55.5$ Relative lowering of VP = x_2 (where x_2 is mole fraction of solute) Relative lowering of VP = n_2/n_1+n_2 (n_2 is no. of moles of solute , n_1 is no. of moles of solvent)	1/2	
	= 2/2+55.5 = 2/57.5 = 0.034	1/2	
19	 (i)t_{2g}⁴e_g² Paramagentic (ii)Dichloridobis(ethane-1,2-diamine)cobalt(III)nitrate OR 	¹ / ₂ , ¹ / ₂ 1	
	(i)Square planar (ii)Cu ²⁺ = $3d^9 1$ unpaired electron so $\sqrt{1(3)} = 1.73BM$	1	
20	Reaction is a complex reaction. Order of reaction is 1.5.	1/2	
	Molecularity cannot be 1.5, it has no meaning for this reaction. The reaction occurs in steps, so it is a complex reaction. (ii)units of k are $mol^{-1/2}L^{1/2}s^{-1}$	1/2 1	

	OR	
	Ans : let the rate law expression be Rate = $k [P]^{x}[Q]^{y}$	
	from the table we know that	
	Rate $1 = 3.0 \times 10^{-4} = k (0.10)^{x} (0.10)^{y}$	
	Rate $2 = 9.0 \times 10^{-4} = k (0.30)^{x} (0.30)^{y}$	
	Rate $3 = 3.0 \times 10^{-4} = k (0.10)^{x} (0.30)^{y}$	
	Rate 1/ Rate $3 = (1/3)^y$ or $1 = (1/3)^y$	
	So $y = 0$	1/2
	Rate 2/ Rate $3 = (3)^x$ or $3 = (3)^x$	/ 2
	So x = 1	1/2
	Rate = $k [P]$	1
21	$k = 0.693/t_{1/2}$ k = 0.693/5730 years ⁻¹	1/2
	$t = 2.303 \log Co$	72
	$\frac{2.300}{k}$ $\frac{100}{Ct}$	1/2
	let Co = 1 Ct = $3/10$ so Co/Ct = $1/(3/10) = 10/3$	
	$t = 2.303 \times 5730 \log \frac{10}{2}$	1/2
	$\begin{array}{c} 0.693 & 3 \\ t = 19042 \text{ x} (1-0.4771) = 9957 \text{ years} \end{array}$	1/2
	t = 19042 x (1-0.4771) = 9937 years	/2
22	$CH_3 - CH - CH_3 \xrightarrow{H^+} CH_3 - CH - CH_3$	
	 CH ₃ OH CH ₃ OH ₂	1/2
	an an the an	1/2
	$\begin{array}{c} CH_3 - CH - CH_3 CH_3 - CH_3 CH_3 - CH_3 CH_3 \xrightarrow$	72
	ĊH ₃ ĆH ₃ ĆH ₃	
	H \downarrow $+$ $ 12 - hydride shift + t$	1/
	$CH_3 - C - CH_3 - CH_3 \xrightarrow{12 - hydride shilt} CH_3 - CH_2 - CH_3$	1/2
	CH ₃ CH ₃	
	Br	
	$CH_3 - \overset{+}{C} - CH_2 - CH_3 \qquad Br^- \longrightarrow CH_3 - \overset{-}{C} - CH_2 - CH_3$	1/2
	CH ₃ CH ₃	
	accume accuments	
23	XeF ₆	1
	Central atom Xe has 8 valence electrons, it forms 6 bonds with F and has	
	1 lone pair. According to VSEPR theory, presence of 6 bp and 1 lp results in	1
	distorted octahedral geometry	-
L		l

	F F F F F F	
24.	Racemic mixture will be given by 2 chlorobutane as it is an optically active compound.	1
	When 2 chlorobutane undergoes S_N^1 reaction, both front and rear attack are possible, resulting in a racemic mixture	1
25	Let no. of Atoms of element P be x No. of tetrahedral voids = 2x No. Of octahedral voids = x	1/2
	Atoms of Q = $1/3 (2x) + x = 5x/3$ P _x Q _{5x/3}	1/2
	P ₃ Q ₅	1
26	SECTION C	
26	 (i)Due to large surface area and ability to show variable oxidation states (ii)Due to high value of third ionisation enthalpy (iii) Mo(VI) and W(VI) are more stable than Cr(VI). OR (i) The general trend towards less negative E° V values across the series is related to the general increase in the sum of the first and second ionisation enthalpies. (ii) The high energy to transform Cu(s) to Cu²⁺ (aq) is not balanced by its hydration enthalpy. 	1 1 1
	(iii) The stability of the half-filled d sub-shell in Mn ²⁺ and the completely filled d^{10} configuration in Zn ²⁺ are related to their more negative E^o V values	1
27	(i) Aniline, <i>N</i> -ethylethanamine, Etanamine (ii)Ethanamine, ethanol, ethanoic acid	1 1
	(iii) N, N dimethylmethanamine, methanamine, N-methylmethanamine OR	1
	(i) N-methyletahnamine is a secondary amine. When it reacts with benzenesulphonyl chloride, it forms N- Ethyl -N methyl sulphonamide while and	1

N,N-dimethyl etahnanmine is a tertiary ami benzenesulphonyl chloride.	N,N-dimethyl etahnanmine is a tertiary amine it does not react with		
(ii) NO ₂ NH ₂	NH ₂		
$\bigcirc \qquad \xrightarrow{H_2/N_1} \qquad \bigcirc \qquad \xrightarrow{Br_2/H_2Q}$	Br Br 1		
(iii)Butan-1-ol Alcohol forms stronger hydrogen bonds wi higher electronegativity of O in alcohol that			
	/2 1/2		
28 We know that $d = zM/N_a a^3$ For fcc, $z=4$ therefore $d = 4$ for bcc, $z=2$ therefore $d' = 2 x M / Na (3.0 d/d' = 4/(3.5 x 10^{-8})^3 / 2/(3.0 x 10^{-8})^3 = 1.2$	1×10^{-8}) ³ g/cm ³		
$\begin{array}{c ccccccccccccccccccccccccccccccccccc$	$\begin{array}{c c} (CH_2)_4 - NH_2 \\ \\ -C - CH - NH_2 \end{array} \qquad 1$		
$\begin{array}{c c} CH_2COOH & CH_3 \\ & \\ HOOC - CH - N - C - CH - N - C \\ & & \\ H & O & H \end{array}$	$- CH - NH_2$		
(ii) $H_{3}^{+}N - C - COO^{-}$ $H_{3}^{+}O - COO^{-}$	1		
30 i. Arrange the following in decreasing ord $Cl_2 > Br_2 > F_2 > I_2$	er of bond dissociation enthalpy 1		
ii. Bi does not form $p\pi$ - $p\pi$ bonds as its effective overlapping is not possible iii.Due to small size of oxygen, it has g	reater electron electron repulsions		
SECT	- 1		
31. (i) (a) $3Cu + 8 HNO_3(dilute) \rightarrow 3C_{(b)}$			





