## MARKING SCHEME

## PRACTICE PAPER - 01

## SECTION A

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

## SECTION B, C, D

Q.No.	VALUE POINTS	MARKS
	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
	$\begin{array}{c} & & & \\ & & & & \\ & & & \\ & & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ &$	
	$= \underbrace{\begin{pmatrix} \Theta \\ Fast step \\ Fast step \\ \Theta \\ $	
	OR	
	(i) $NH_2$ $NaNO_2 + HCl$ $273 - 278 K$ $Cl$ $Cu_2Cl_2$	1
	(ii) CH <sub>3</sub> CH(Br)CH <sub>3</sub> alc KOH CH <sub>3</sub> CH=CH <sub>2</sub> HBr, organic peroxide CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> Br	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. 2 m ag solution of urea means 2 moles of urea in 1kg of water.	1
	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 <sub>2g</sub> <sup>4</sup> eg <sup>2</sup> Paramagentic (ii)Dichloridobis(ethane-1,2-diamine)cobalt(III)nitrate OR	<sup>1</sup> / <sub>2</sub> , <sup>1</sup> / <sub>2</sub> 1
	(i)Square planar (ii)Cu <sup>2+</sup> = 3d <sup>9</sup> 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

	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 $3 = (3)^x$	
	So $x = 1$	1/2
	Rate = k [P]	1
21	$\mathbf{k} = 0.693/t_{1/2}$	
	$k = 0.693/5730 \text{ years}^{-1}$	1/2
	$t = 2.303 \log CO$	1/2
	k Ct let Co = 1 Ct = $3/10$ so Co/Ct = $1/(3/10) = 10/3$	
	$t = 2.303 \times 5730 \log 10$	1/2
	0.693 3	
	$t = 19042 \times (1-0.4771) = 9957$ years	1/2
22	$CH_3 - CH - CH - CH_3 \xrightarrow{H^*} CH_3 - CH - CH - CH_3$	
	 СН3 ОН СН3 ОН2	1/2
		1/2
	$\begin{array}{c} CH_3 - CH - CH_3 - CH_3  CH_3 - CH_3  CH_3 \xrightarrow$	12
	CH, CH, CH,	
	$CH_1 - C - CH_2 - CH_3 \xrightarrow{12-hydride shift} CH_3 - CH_2 - CH_3$	1/2
	 CH <sub>3</sub> CH <sub>3</sub>	
	Br	12
	$CH_3 - C - CH_2 - CH_3 = CH_3 - C - CH_2 - CH_3$	1/2
	CH <sub>3</sub> CH <sub>3</sub>	
23	XeF <sub>6</sub>	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	
	distorted octahedral geometry	1
	uisioned octanedral geometry	

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 = $2x$	1/2
	No. Of octahedral voids = $x$	
	Atoms of $Q = 1/3 (2x) + x = 5x/3$	1/2
	P <sub>x</sub> Q <sub>5x/3</sub>	
	P <sub>3</sub> Q <sub>5</sub>	1
26	SECTION C	
20	(i)Due to large surface area and ability to show variable oxidation states	1
	(ii)Due to high value of third ionisation enthalpy	1
	(iii) Mo(VI) and W(VI) are more stable than Cr(VI).	1
	OR	
	(i) The general trend towards less negative E <sup>o</sup> V values across the series is related to the general increase in the sum of the first and second	
	ionisation enthalpies.	1
	<li>(ii) The high energy to transform Cu(s) to Cu<sup>2+</sup> (aq) is not balanced by its</li>	•
	hydration enthalpy.	
	(iii) The stability of the half-filled d sub-shell in Mn <sup>2+</sup> and the completely filled d <sup>10</sup> configuration in Zn <sup>2+</sup> are related to their more negative E <sup>o</sup> V values	1
	Tulues	1
27	(i) Aniline, N-ethylethanamine, Etanamine	1
	(ii)Ethanamine, ethanol, ethanoic acid	1
	(iii) N, N dimethylmethanamine, methanamine, N-methylmethanamine OR	1
	(i) N-methyletahnamine is a secondary amine. When it reacts with	1
	benzenesulphonyl chloride, it forms N- Ethyl -N methyl sulphonamide while and	

	N,N-dimethyl etahnanmine is a tertiary amine it does not react with	
	benzenesulphonyl chloride.	
	(ii) NO <sub>2</sub> NH <sub>2</sub> NH <sub>2</sub>	
	$\bigcirc \qquad \xrightarrow{H_2/N_1} \bigcirc \qquad \xrightarrow{Br_2/H_2O} \bigcirc \bigcirc$	
		1
	(iii)Butan-1-ol	
	Alcohol forms stronger hydrogen bonds with water than formed by amine due to	
	higher electronegativity of O in alcohol than N in amine	1.1 m
		1/2
		1/2
		12
28	We know that $d = zM/N_a a^3$ For fcc, z=4 therefore $d = 4 \times M / Na (3.5 \times 10^{-8})^3 \text{ g/cm}^3$	1/2
	For bcc, $z=2$ therefore d' = 2 x M / Na (3.0 x 10 <sup>-8</sup> ) <sup>3</sup> g/cm <sup>3</sup>	
20	$d/d' = 4/(3.5 \times 10^{-8})^3 / 2/(3.0 \times 10^{-8})^3 = 1.26:1$	1/2
29	(i) $CH_3$ $CH_2$ -COOH $(CH_2)_4$ - $NH_2$	1
	HOOC $-CH - N - C - CH - N - C - CH - NH_2$	1
	н о н о	
	CH <sub>2</sub> COOH CH <sub>3</sub> (CH <sub>2</sub> ) <sub>4</sub> - NH <sub>2</sub>	1
	HOOC $-CH - N - C - CH - N - C - CH - NH_2$	
	но но	1
	(ii)	1
	H +	
	$H_3N-C-COO^{-1}$	
	CH <sub>3</sub>	
30	i. Arrange the following in decreasing order of bond dissociation enthalpy	1
	$Cl_2 > Br_2 > F_2 > I_2$	
	ii. Bi does not form $p\pi$ - $p\pi$ bonds as its atomic orbitals are large and diffuse so	1
	effective overlapping is not possible iii.Due to small size of oxygen, it has greater electron electron repulsions	1
	SECTION D	1
31.	(i)	
	(a) $3Cu + 8 HNO_3(dilute) \rightarrow 3Cu(NO_3)_2 + 2NO + 4H_2O$	1
	(b)	





