

## <u>Chemistry Marking scheme</u> <u>Delhi - 2016</u> <u>Set – 56/1/3/D</u>

	<u>Set - 30/1/3/D</u>			
Q.No	VALUE POINTS	MARKS		
1	Ferromagnetism			
2	$CH_3CH_2CH(Cl)CH_3$ ; secondary halide/ 2 <sup>0</sup> carbocation is more	1/2, 1/2		
	stable			
3	NH <sub>3</sub>	1		
4	Like Charged particles cause repulsion/ Brownian motion/ solvation	1		
5	2,4,6-Tribromoaniline / 2,4,6-Tribromobenzenamine	1		
6	(i) $[Cr(H_2O)_5Cl]Cl_2.H_2O$	1		
	<ul> <li>(ii) pentaaquachloridoChromium(III) chloride monohydrate (or chloride hydrate)</li> </ul>	1		
	(no deduction for not writing hydrate)			
7	(i) Mercury cell	1/2		
	(ii) Fuel cell	1/2		
	(iii) Lead storage battery	1/2		
	(iv)Dry cell	1/2		
8	A-Na <sub>2</sub> CrO <sub>4</sub>	1/2		
	$B-Na_2Cr_2O_7$	1/2		
	$C-K_2Cr_2O_7$	1/2		
	Use- strong oxidising agent / as a primary standard in	1/2		
	volumetric analysis			
	OR			
8	$8MnO_4^- + 3S_2O_3^{2-} + H_2O \longrightarrow 8MnO_2 + 6SO_4^{2-} + 2OH^-$	1		
	$Cr_2O_7^{2-}$ + 14 H <sup>+</sup> + 3 Sn <sup>2+</sup> $\rightarrow$ 2 Cr <sup>3+</sup> + 3 Sn <sup>4+</sup> + 7 H <sub>2</sub> O	1		

9	(i) $CH_3-CH_2-\ddot{O}-H + H' \rightarrow CH_3-CH_2-\ddot{O}-H$	1⁄2
	(ii) $CH_3CH_2 = \overset{\bigcirc}{O} + CH_3 = CH_2 = \overset{\frown}{O} + \overset{H}{H} \rightarrow CH_3CH_2 = \overset{\frown}{O} - CH_2CH_3 + H_2O$	1
	(iii) $CH_{3}CH_{3} \xrightarrow{\bullet}_{i} - CH_{3}CH_{3} \longrightarrow CH_{3}CH_{3} \xrightarrow{\bullet}_{i} - CH_{2}CH_{3} + H^{*}$	1/2
10	(i) zero order, bimolecular/ unimolecular	1/2, 1/2
	(ii) mol L <sup>-1</sup> s <sup>-1</sup>	1
11	$\log k = \log A - E_a/2.303RT$	1/2
	$E_a / 2.303 \text{ RT} = 1.0 \times 10^4 \text{ K/ T}$	
	$E_a = 1.0 \times 10^4 \times 2.303 \times 8.314$	
	=191471.4 J/mol	1
		1000
	$t_{1/2} = 0.693 / k$	1/2
	k = 0.693/200  min	
	$= 0.0034 \text{min}^{-1}$	1



	• (i)	_
	Adsorption Absorption	
	Surface phenomena Bulk phenomena	
	The accumulation ofThe substance is uniformly	
	molecular species distributed throughout	
	at the surface rather than in the bulk of the solid	
	the bulk of a solid or liquid is essentially a bulk	
	termed adsorption. phenomenon. (any one difference)	1
	(any one difference)	_   `
	(ii) AlCl <sub>3</sub> , more positive charge/Hardy-Schulze rule	$\frac{1}{2} + \frac{1}{2}$
	(iii)Sulphur	1
13	(i) In chlorobenzene, each carbon atom is sp <sup>2</sup> hybridised /	1
	resonating structures / partial double bond character.	
	(ii) Due to +R effect in chlorobenzene/ difference in	1
	hybridization i.e. sp <sup>2</sup> and sp <sup>3</sup> respectively/ -I and +R effect	
	oppose each other while -I effect is the only contributing	
	factor in cyclohexane.	
*	(iii)Due to formation of planar carbocation/ Carbon in	1
	carbocation formed is $sp^2$ hybridised.	
14	$2 \times 10^{24}$ atoms weigh = 300g	1 .
	$6.022 \times 10^{23}$ atoms weigh = $(300 \times 6.022 \times 10^{23})/2 \times 10^{24}$	
	= 90.3 g	
	d = z x M	
	$d = \frac{z \times M}{a^3 N_A}$	$\frac{1}{2} + \frac{1}{2}$
	103	
	$= 4 \times 90.3 / (250 \times 10^{-10} \text{ Jx} \text{ N}_0)$	/2 . /2
	$= 4x90.3/(250x10^{-10})xN_0$ =38.4 gcm <sup>-3</sup>	
	$= 4x90.3/(250x10^{-10})XN_0$ =38.4 gcm <sup>-3</sup> (or any other correct method	1
	$= 4x90.3/(250x10^{-10})xN_0$ =38.4 gcm <sup>-3</sup> (or any other correct method	1
15	(or any other correct method	) 1
15	$= 4x90.3/(250x10^{-10})xN_0$ =38.4 gcm <sup>-3</sup> (or any other correct method (i) ability of oxygen to form multiple bond/ pπ-dπ bond. (ii) Partially filled d orbitals / due to comparable energies of ns	1
15	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> </ul>	) 1
15	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of</li> </ul>	) 1 1 1 1 1
÷.	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> </ul>	) 1 1 1 1 1
ž.	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> <li>(i) CH<sub>3</sub>OH, (CH<sub>3</sub>)<sub>3</sub>C-I</li> </ul>	) 1 1 1 1 1
÷.	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> </ul>	1 1 1 1 1 1 1 1
÷.	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> <li>(i) CH<sub>3</sub>OH, (CH<sub>3</sub>)<sub>3</sub>C-I</li> <li>(ii) CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>OH</li> </ul>	1 1 1 1 1 1
÷.	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> <li>(i) CH<sub>3</sub>OH, (CH<sub>3</sub>)<sub>3</sub>C-I</li> <li>(ii) CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>OH</li> </ul>	1 1 1 1 1 1 1 1
16	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> <li>(i) CH<sub>3</sub>OH, (CH<sub>3</sub>)<sub>3</sub>C-I</li> <li>(ii) CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>OH</li> </ul>	1 1 1 1 1 1 1 1 1
16	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> <li>(i) CH<sub>3</sub>OH, (CH<sub>3</sub>)<sub>3</sub>C-I</li> <li>(ii) CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>OH</li> <li>(iii) CH<sub>3</sub> CH<sub>2</sub>CH<sub>2</sub>OH</li> <li>(ii) Zone refining</li> </ul>	1 1 1 1 1 1 1 1 1 1
16	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> <li>(i) CH<sub>3</sub>OH, (CH<sub>3</sub>)<sub>3</sub>C-I</li> <li>(ii) CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>OH</li> <li>(iii) Cone refining</li> <li>(ii) Leaching / Bayer's process</li> </ul>	1 1 1 1 1 1 1 1 1 1 1 1 1 1
16	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> <li>(i) CH<sub>3</sub>OH, (CH<sub>3</sub>)<sub>3</sub>C-I</li> <li>(ii) CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>OH</li> <li>(iii) CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>OH</li> <li>(iii) CHaching / Bayer's process</li> <li>(iii) Reducing agent / to form CO which acts as a reducing agent.</li> </ul>	1 1 1 1 1 1 1 1 1 1
16	<ul> <li>(or any other correct method</li> <li>(i) ability of oxygen to form multiple bond/ pπ-dπ bond.</li> <li>(ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals</li> <li>(iii) due to relative stabilities of the f<sup>0</sup>, f<sup>7</sup> and f<sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals.</li> <li>(i) CH<sub>3</sub>OH, (CH<sub>3</sub>)<sub>3</sub>C-I</li> <li>(ii) CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>OH</li> <li>(iii) CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>OH</li> <li>(iii) CHaching / Bayer's process</li> <li>(iii) Reducing agent / to form CO which acts as a reducing agent.</li> </ul>	1 1 1 1 1 1 1 1 1 1 1 1 1 1
16	(or any other correct method (i) ability of oxygen to form multiple bond/ $p\pi$ -d $\pi$ bond. (ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals (iii) due to relative stabilities of the f <sup>0</sup> , f <sup>7</sup> and f <sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals. (i) CH <sub>3</sub> OH, (CH <sub>3</sub> ) <sub>3</sub> C-I (ii) CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> OH (ii) CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> OH (ii) Leaching / Bayer's process (iii) Reducing agent / to form CO which acts as a reducing agent. (i) E <sup>0</sup> <sub>cell</sub> = E <sup>0</sup> <sub>c</sub> -E <sup>0</sup> <sub>a</sub> -	1 1 1 1 1 1 1 1 1 1 1 1 1 1
16	(or any other correct method (i) ability of oxygen to form multiple bond/ $p\pi$ -d $\pi$ bond. (ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals (iii) due to relative stabilities of the f <sup>0</sup> , f <sup>7</sup> and f <sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals. (i) CH <sub>3</sub> OH, (CH <sub>3</sub> ) <sub>3</sub> C-I (ii) CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> OH (ii) CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> OH (ii) Zone refining (ii) Leaching / Bayer's process (iii) Reducing agent / to form CO which acts as a reducing agent. (i) E <sup>0</sup> <sub>cell</sub> = E <sup>0</sup> <sub>c</sub> -E <sup>0</sup> <sub>a</sub> =(-0.44)-(-0.74) V	1 1 1 1 1 1 1 1 1 1 1 1 1 1
16	(or any other correct method (i) ability of oxygen to form multiple bond/ $p\pi$ -d $\pi$ bond. (ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals (iii) due to relative stabilities of the f <sup>0</sup> , f <sup>7</sup> and f <sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals. (i) CH <sub>3</sub> OH, (CH <sub>3</sub> ) <sub>3</sub> C-I (ii) CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> OH (ii) CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> OH (ii) Leaching / Bayer's process (iii) Reducing agent / to form CO which acts as a reducing agent. (i) E <sup>0</sup> <sub>cell</sub> = E <sup>0</sup> <sub>c</sub> -E <sup>0</sup> <sub>a</sub> -	1 1 1 1 1 1 1 1 1 1 1 1 1 1
16	(or any other correct method (i) ability of oxygen to form multiple bond/ $p\pi$ -d $\pi$ bond. (ii) Partially filled d orbitals / due to comparable energies of ns and (n-1) d orbitals (iii) due to relative stabilities of the f <sup>0</sup> , f <sup>7</sup> and f <sup>14</sup> occupancies of the 5f orbitals/ Comparable energies of 7s,6d,5f orbitals. (i) CH <sub>3</sub> OH, (CH <sub>3</sub> ) <sub>3</sub> C-I (ii) CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> OH (ii) CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> OH (ii) Zone refining (ii) Leaching / Bayer's process (iii) Reducing agent / to form CO which acts as a reducing agent. (i) E <sup>0</sup> <sub>cell</sub> = E <sup>0</sup> <sub>c</sub> -E <sup>0</sup> <sub>a</sub> =(-0.44)-(-0.74) V	1 1 1 1 1 1 1 1 1 1 1 1 1 1



	$E_{cell} = E_{cell}^0 - 0.059 \log [Cr^{3+}]^2$ n [Fe <sup>3+</sup> ] <sup>3</sup>	1/2
	ii [ie]	1
	$F_{\mu\nu} = F^0_{\mu\nu} = 0.059 \log [0.01]^2$	
	$E_{cell} = E_{cell}^0 - 0.059 \log [0.01]^2$ 6 [0.1] <sup>3</sup>	
	= 0.30 - (-0.059/6)	1
	=0.3098V	
19	(i) $\beta$ -D glucose and $\beta$ -D-galactose / glucose and galactose	16 16
	(ii) water soluble ,excreted out of the body	$\frac{1}{2}, \frac{1}{2}$
	(iii)In nucleotide, phosphoric acid/phosphate group attached to	1
	the nucleoside / structures of both nucleoside and nucleoside /	
	nucleotide= base +sugar + phosphate group, nucleoside= base	1
	+sugar.	
20	d <sup>2</sup> sp <sup>3</sup> , Paramagnetic, low spin	1.1/.1/
20		$1, \frac{1}{2}, \frac{1}{2}$
		1
	en pi	1
•	( en'	
21	(i) $C_6H_5NH_2$ , $C_6H_5N_2^+CI^-$ , $C_6H_5I$	$\frac{1}{2} + \frac{1}{2} + \frac{1}{2}$
	(ii) CH <sub>3</sub> CN, CH <sub>3</sub> CH <sub>2</sub> NH <sub>2</sub> , CH <sub>3</sub> CH <sub>2</sub> NC	$\frac{1}{2} + \frac{1}{2} + \frac{1}{2}$
22	a. Catalyst / initiator of free radical	1
	b. Hexamethylene diamine and adipic acid / structure /	1/2, 1/2
	IUPAC name	,
	c. Buna-S <polystyrene<terylene< td=""><td>1</td></polystyrene<terylene<>	1
	OR	
22	Chain initiation steps	
22		
	$C_{H_{a}} \xrightarrow{C_{c}} O'O \xrightarrow{C_{c}} C_{c}H_{a} \xrightarrow{C_{c}} 2C_{c}H_{a} \xrightarrow{C_{c}} O \xrightarrow{O} \xrightarrow{C_{c}} 2\hat{C}_{a}H_{a}$ Berzeyl peroxide Phenyl radical	1
	$C_{H_1}$ +CH <sub>2</sub> =CH <sub>2</sub> $\longrightarrow$ C <sub>4</sub> H <sub>4</sub> -CH <sub>2</sub> -CH <sub>2</sub>	
	Chain propagating step	
	$C_{4}H_{3} - CH_{2} - CH_{3} + CH_{3} = CH_{7} \longrightarrow C_{4}H_{3} - CH_{3} - CH$	
		1
	$\mathbf{C}_{\mathbf{s}}\mathbf{H}_{\mathbf{s}} + \mathbf{C}\mathbf{H}_{\mathbf{s}} - \mathbf{C}\mathbf{H}_{\mathbf{s}} + \mathbf{C}\mathbf{H}_{\mathbf{s}} - \mathbf{C}\mathbf{H}_{\mathbf{s}} + \mathbf{C}\mathbf{H}_{\mathbf{s}} - \mathbf{C}\mathbf{H}_{\mathbf{s}} + C$	1
	Chain terminating step	
	For termination of the long chain, these free radicals can combine in different ways to form polythene. One mode of termination of chain is shown as under:	
	in different ways to form polythene. One mode of termination of	1

23	(i)Aware, concerned or any other correct two values.	$\frac{1}{2} + \frac{1}{2}$
	(ii) Side effects, unknown health problems	1
	(iii) Neurologically active drugs/ stress relievers	1
	Example- valium, equanil	$\frac{1}{2} + \frac{1}{2}$
	(or any other correct two example)	
24	A -C <sub>6</sub> H <sub>5</sub> COCH <sub>3</sub>	1
	$B-C_6H_5CH_2CH_3$	1
	C-C <sub>6</sub> H <sub>5</sub> COOH	1



	D ,E -C <sub>6</sub> H <sub>5</sub> COONa , CHI <sub>3</sub>	1+1
	OR	
24	a)HCHO + HCHO conc MACH HCOONa +CH <sub>3</sub> OH	1
	(or any other example)	
	b)CH <sub>3</sub> CH=N-NHCONH <sub>2</sub>	1
	c) Stronger -I effect of fluorine ,stronger acid less pka / strong	1
	electron withdrawing power of fluorine.	
	d)CH <sub>3</sub> CH=CHCH <sub>2</sub> CHO	1
	e)Silver mirror formed on adding ammonical silver nitrate to	1
	propanal and not with propanone (or any other correct test)	

25	a) $\Delta T_{f} = i \frac{K_{f} w_{b} x1000}{K_{f} w_{b} x1000}$	1
	M <sub>b</sub> x w <sub>a</sub>	
		1
	$\Delta T_{f} = 3 \times (1.86 \times 1.9/95 \times 50) \times 1000$	
	= 2.23K	
	$T_{f} - \Delta T_{f} = 273.15 - 2.23 / 273 - 2.23$	1
	$T_{f} = 270.92 \text{ K or } 270.77 \text{ K}$	1
	b)	NUMBER AND ADDRESS
	i)2M glucose ; More Number of particles / less vapour pressure	$\frac{1}{2} + \frac{1}{2}$
	ii)Reverse Osmosis	1
	OR	
	a) AT = 16 m m1000	
	$\Delta T_{f} = \frac{K_{f} w_{b} \times 1000}{M_{b} \times w_{a}}$	1
		1
	$0.383 = (3.83 \times 2.56/M \times 100) \times 1000$	1
	M=256	
	$S \times x = 256$	
	$32 \times x = 256$	
	x=8	1
	b)	1
	i)Shrinks	1
	ii)swells	1

26	<u>a)</u>		
	i.	Endothermic compound / decomposition of ozone is exothermic	1
		in nature and $\Delta { extsf{G}}$ is negative / decomposition of ozone is	
		spontaneous.	





