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# **AIEEE 2007 Chemistry Answers and Solutions**

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#### **Some Important Hints and Solutions:**

1)4	2) 3	3) 3	4) 1
5)4	6) 4	7) 1	8) 1
9)1	10) 4	11) 3	12) 4
13) 2	14) 4	15) 3	16) 4
17) 3	18) 2	19) 4	20) 1
21) 2	22) 3	23) 2	24) 2
25) 4	26) 4	27) 1	28) 1
29) 1	30) 3	31) 3	32) 4
33) 1	34) 3	35) 4	36) 3
37) 3	38) 3	39) 1	40) 4

#### **Some Important Hints and Solutions:**

- **1**  $\Delta H = E_f E_b$
- $\Delta H = 80-100 = -20$
- $nFE^{\circ}_{cell} = -2.303 \text{ RT} \log(Zn^{2+}/Cu)$ 2 98×log(Zn<sup>2+</sup>/Cu<sup>2+</sup>) => 2×96500×1.10 = 2.303×8.3  $=> 37.3 = \log(Zn^{2+}/Cu^{2+})$  $=> \log(Zn^{2+}/Cu^{2+}) = 10^{37}$ 3 Henderson-Hasselbalch equation:  $pH = pK_a + log10$  [Conjugate Base]/[Acid] 5c/0.5c

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pH = 4.5

 $pH = 4.5 + \Omega q^2$ 

=9.5



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**4** A first order reaction has a rate proportional to the concentration of one of reactants.

In this case the rate of reaction is dependent on A and not dependent of So this is first order reaction.

So Rate = k[A][B]

= mole/(liter.sec) = k(mole/liter)<sup>2</sup>

= k = mole<sup>-1</sup>liter.sec<sup>-1</sup>

**5** 4f is shielded more than 5f

**6** Rh(I), Ir(I), Pd(II), Pt(II), and Au(III) belongs to square planar geometry.

7 The asymmetric molecules that have no center, axis of symmetry rotates the plane of polarized light.

The simplest type of asymmetric molecules (chirat holecule) is one which have four different group attached to same carbon item.

Compound (1) does not have any plane of symmetry, so it is optically active.

**8** Secondary structure in proteins confists a helics and  $\beta$ -sheets structures.

These structures are formed as a result of H-bonding between different peptide groups.

of H  $CH_3 - C \equiv CH + HBr$ C-CH, Br 9 Rr

**10** This is a Carbylan necesction. The Carbylamine reaction is a chemical test for detection of primary amines So the outputs are:  $C_2H_5NC$  and 3KCl

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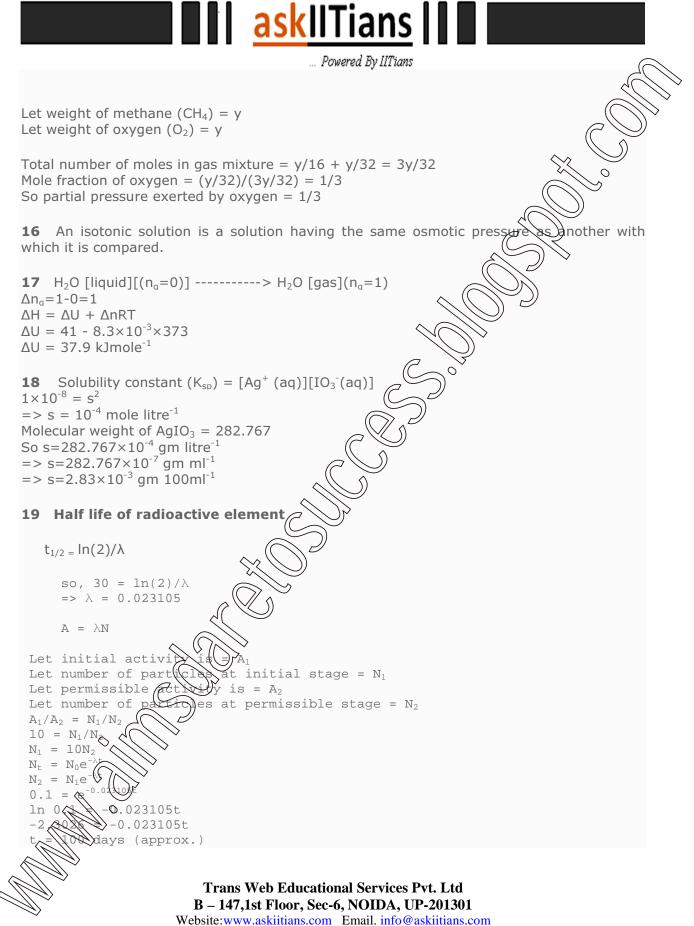
## **ask**IITians ... Powered By IITians 11 $CH_3$ $CH_3$ $CH_3$ **12** If a group takes a negative charge by induction, the group possess a -I effect. If a group takes a positive charge by induction, the group possess a +I effect. If a group takes a negative charge by resonance, the group possess a -R effect. If a group takes a positive charge by resonance, the group 463333 +R effect. $-NO_2$ group in benzene ring shows -I and -R effect, which deactivates the ring towards electrophilic substitution but activates it towards nucleophilic substitution. **13** Bond order is the number of bonds between а atoms. pair of Bond order= 1/2(number of bonding orbital - number of)anti bonding orbital) In C<sub>2</sub>, bond order=2 (diamagnetic) In $C_2^+$ , bond order=1.5 (paramagnetic) In NO, bond order=2.5 (paramagnetic) In NO<sup>+</sup>, bond order=3 (diamagnetic In $O_2$ , bond order=2 (diamagnetic In $O_2^+$ , bond order=2.5 (parama In N<sub>2</sub>, bond order=3 (diamagnetic) In $N_2^+$ , bond order=2.5 (paramagnetic)

**14** The actinoids exhibit more number of oxidation states than lanthanoids because 5f orbitals is further from the nucleus than the 4f orbitals. So, 5f orbital electrons are held less strongly than the 4f orbital electrons.

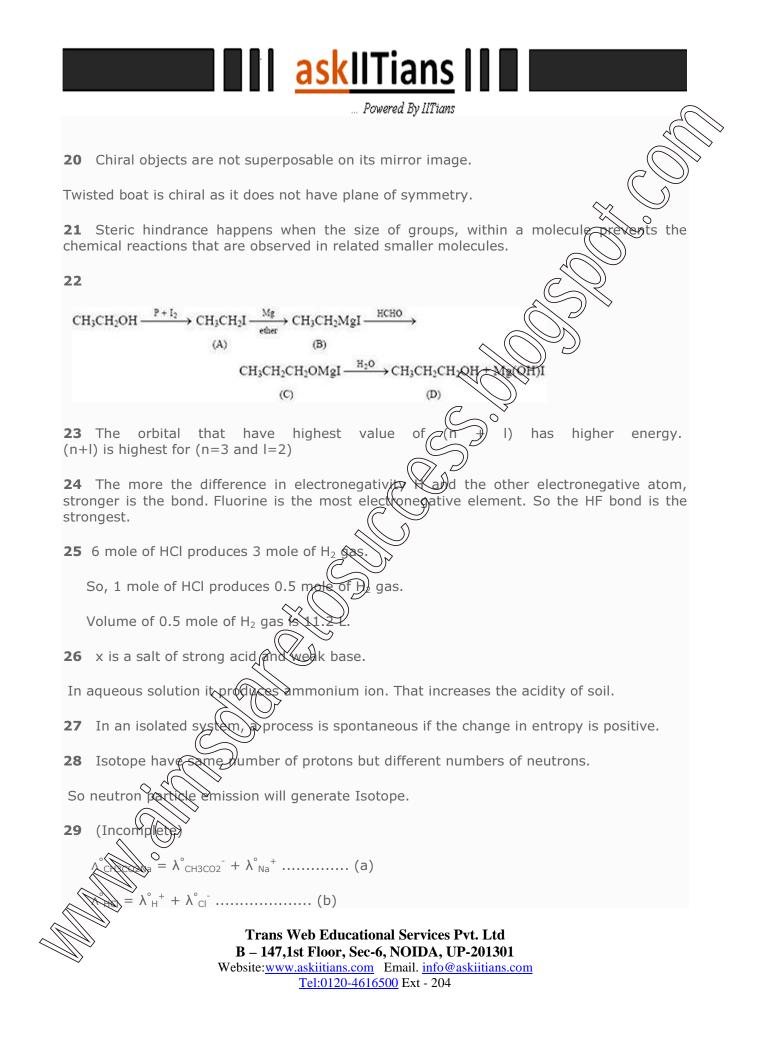
**15** In a most the orideal gases, partial pressure is the pressure which the gas would have if it would have alone occupied the volume. The total pressure of a gas mixture is the sum of the partial pressures of each individual gas in the mixture.

Molecular weight of methane  $(CH_4) = 16$ Molecular weight of oxygen  $(O_2) = 32$ 

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Add (a) and (b)  $\Lambda^{\circ}_{CH3CO2Na} + \Lambda^{\circ}_{HCI} = \lambda^{\circ}_{CH3CO2} + \lambda^{\circ}_{Na} + \lambda^{\circ}_{H} + \lambda^{\circ}_{CI}$   $=> \Lambda^{\circ}_{CH3CO2Na} + \Lambda^{\circ}_{HCI} = \lambda^{\circ}_{CH3CO2} + \lambda^{\circ}_{NaCI} + \lambda^{\circ}_{H}^{+}$  $=> \Lambda^{\circ}_{CH3CO2H} = \lambda^{\circ}_{CH3CO2} + \lambda^{\circ}_{H}^{+}$ 

**30** overall basic strength varies as  $2^{\circ} > 1^{\circ} > 3^{\circ}$ 

**31** In aliphatic compounds, carbon atoms are joined together in straight enains, branched chains, or non-aromatic rings

Any aliphatic carbon with hydrogen attached to it, in combination with benzene ring, will be oxidized to benzoic acid by  $KMnO_4/H^+$ 

32 2 5 3-ethyl-4, 4-dimethylheptane s)σ2P<sub>z</sub><sup>2</sup> , π2P<sub>x</sub><sup>2</sup> **33**  $O_2^{2^-} = 18 \sigma 1s^2 \sigma^* 1s^2 \sigma 2s^2$  $= \pi 2 P_v^2$ ,  $\pi^* 2 P_x^2 = \pi^* 2 P_v^2$ Hence diamagnetic 34 Due to inert pair effect +2 oxidation state increases as we move down this group in group 14.  $\propto$  SnX<sub>2</sub> << PbX<sub>2</sub> So, SiX<sub>2</sub>  $3B_{R_{1}} + 6M_{2}OH \rightarrow 5NaBr + NaBrO_{3} + 3H_{2}O$ 35 **Trans Web Educational Services Pvt. Ltd** B – 147,1st Floor, Sec-6, NOIDA, UP-201301 Website:www.askiitians.com Email. info@askiitians.com Tel:0120-4616500 Ext - 204

# *... Powered By IITians* **36** Greater the charge/size ratio of a cation, the mote is its polarizing power.

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So 3 is correct

**37** Molecular weight of  $H_2SO_4 = 98$ Molarity of solution given = 3.6 So 1 liter of solution contains 3.6 moles of  $H_2SO_4$ So 1 liter of solution contains  $3.6 \times 98$  gm of  $H_2SO_4$ Let the density of solution = D gm/ml So 1000D gm of solution contains  $3.6 \times 98$  gm of  $H_2SO_4$ Now  $(3.6 \times 98)/(1000D) = 0.29$ So D = 1.2 gm/ml

**38**  $K = K_1 \times K_1$ 

 $= K = 1.0 \times 10^{-5} \times 5.0 \times 10^{-10}$ 

 $= > K = 1.0 \times 10^{-15}$ 

**39** As per Raoult's law:

Vapor pressure of an ideal solution is dependent on the vapor pressure of each chemical component and the mole fraction of the component present in the solution. So,  $290 = 200 \times 0.4 + P \times 0.6$ 

**40** 
$$\Delta G = \Delta H - T\Delta S = 0$$

$$=> \Delta H = T\Delta S$$

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