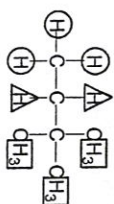
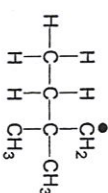
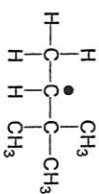
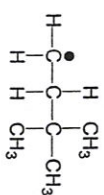


1	C	2	A
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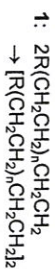
1 C



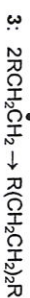
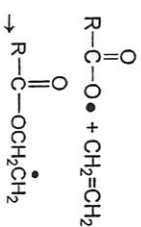
The 3 different forms of X• are:



2 A



2:



3(a) No formation of oxides of sulfur which causes acid rain OR

Efficient combustion prevents photochemical smog (less unburnt hydrocarbon OR less CO being released)

(b) (i) $W_a = 8(1 - 0.89)^3 = 0.0428$ Mass of $\text{C}_8\text{H}_{18} = 0.0428 \times 100 = 4.28 \text{ g}$ M_r of $\text{C}_8\text{H}_{18} = (8 \times 12) + 18 = 114$ Number of moles of $\text{C}_8\text{H}_{18} = 4.28 + 114 = 0.0375 \text{ mol}$

(ii) Isomer Y.

Presence of plane of symmetry in the molecule
Hence rotation at each chiral centre cancels out / mirror images superimposable / no enantiomer

(iii)

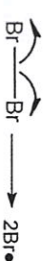
Structural formula	No of hydrogen	Mole ratio
$\begin{array}{c} \text{CH}_3 \quad \text{CH}_3 \\ \quad \\ \text{C}-\text{CH}_2-\text{CH}_2-\text{CH}-\text{CH}-\text{CH}_2-\text{CH}_3 \\ \quad \\ \text{Cl} \quad \text{CH}_3 \quad \text{CH}_3 \end{array}$	6	3
$\begin{array}{c} \text{CH}_3-\text{CH}-\text{CH}-\text{CH}-\text{CH}_2-\text{CH}_3 \\ \quad \quad \\ \text{Cl} \quad \text{CH}_3 \quad \text{CH}_3 \end{array}$	4	2
$\begin{array}{c} \text{CH}_3-\text{CH}_2-\text{C}-\text{CH}-\text{CH}_2-\text{CH}_3 \\ \quad \\ \text{Cl} \quad \text{CH}_3 \end{array}$	2	1
$\begin{array}{c} \text{CH}_3-\text{CH}_2-\text{CH}-\text{CH}-\text{CH}_2-\text{CH}_3 \\ \quad \\ \text{CH}_2\text{Cl} \quad \text{CH}_3 \end{array}$	6	3

4(a) (i) Condition: high temperature or uv light

Free Radical Substitution

Mechanism:

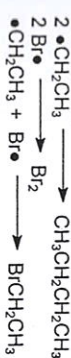
Initiation



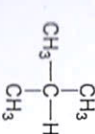
Propagation



Termination



(ii) Alkane B:

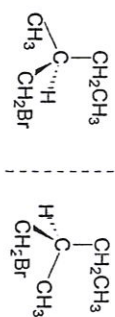
(iii) $\text{C/D} = 2.3 : 1$ or $7 : 3$ or $21 : 9$

Relative rate suggests 21 : 1, but there are 1 tertiary to 9 primary hydrogen atoms, hence this ratio becomes 21 : 9. [1]

[1]

(b) (i) Optical isomerism

[1]



[2]

(ii) I. Boiling point / melting point / density

II. Optical activity / Direction of rotation of plane-polarised light

[1]

[1]

(iii) 2,2-dimethylpropane is more highly branched, thus it is more compact / spherical compared to 2-methylbutane.

[1]

As a result, 2,2-dimethylpropane has a smaller surface area for instantaneous dipole-induced dipole attractions (or Van Der Waals forces) between molecules to operate. Hence, it has a much lower boiling point than 2-methylbutane.

[1]