

15. Halogen compounds

15.1 Halogenoalkanes

Paper 2

Marking Scheme

Q1.

| | | |
|---|--|----------|
| (b) | | 3 |
| <p>M1 correct dipole ($\delta^+C-Br^{\delta-}$) AND curly arrow from bond of C—Br to Br (or just beyond)</p> <p>M2 correct intermediate ALLOW R for $-(CH_2)_5CH_3$</p> <p>M3 curly arrow from lone pair on OH^- to the C of intermediate (or suitable position where C—OH bond is made)</p> <p><i>If SN2 mechanism shown:</i></p> <p>Apply SN2 mechanism mark scheme only if mechanism shows OH^- / NaOH species attacking the reactant and no OH^-</p> <p>M1 correct dipole ($\delta^+C-Br^{\delta-}$) AND curly arrow from bond of C—Br to Br of reactant (or just beyond)</p> <p>NOT M2</p> <p>M3 curly arrow from lone pair on $:OH^-$ to the appropriate C of the reactant</p> | | |

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|---------|------------------------------------|----------|
| (d)(i) | elimination | 1 |
| (d)(ii) | NaOH / sodium hydroxide in ethanol | 1 |

Q2.

| | | |
|----------|--|----------|
| (c)(iii) | $CH_3COCH_3 + 3NaClO \rightarrow 1CHCl_3 + 1CH_3COONa + 2NaOH$ | 1 |
| (c)(iv) | <p>M1 white precipitate/white ppt forms</p> <p>M2 $CHCl_3$ is hydrolysed / chloride ions are released</p> <p>OR</p> <p>Cl^- / chloride ions which are released react with Ag^+</p> <p>OR</p> <p>silver chloride / $AgCl$ is formed</p> | 1 |

Q3.

| | | |
|--------|-------------|----------|
| (c)(i) | | 1 |
| | bromoethene | 1 |

Q4.

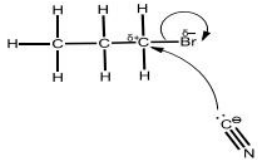
| | | |
|-----|---|----------|
| (e) | $CHCl_3 + 2HF \rightarrow CHCl_2F + 2HCl$ | 1 |
|-----|---|----------|

Q5.

| | | |
|--------|-----------------------|---|
| (a)(i) | KCN in ethanol + heat | 1 |
|--------|-----------------------|---|

Q6.

| | | | | |
|--|----------|---|------------------|---|
| (a) | reaction | reagent and conditions | type of reaction | 6 |
| | 1 | NaOH(aq) [M1] | substitution | |
| | 2 | NH ₃ [M2] in ethanol AND heat under pressure [M3] | substitution | |
| | 3 | NaOH in ethanol AND heat [M4] | elimination | |
| <p>M5 and M6 types of reaction [3 correct types = 2 marks and 2 correct types = 1 mark] All reagents can be identified as (correct) formula or in words</p> | | | | |

| | | |
|--------|--|---|
| (b) | <p>M1 C-I (covalent bond) is weaker or a</p> <p>M2 lower activation energy / lower E_a (with 2-iodopropane)</p> <p>OR</p> <p>M2 explain in terms of the S_N1 mechanism (that is dominant / preferred / occurring with 2-iodopropane is fast(er)) (identical) carbocation is made (more) quickly (with 2-iodopropane)</p> <p>OR</p> <p>low / less energy is required to make the (same) carbocation / intermediate (from 2-iodopropane)</p> <p>OR</p> <p>M2 iodine / I (of C-I) has weaker attraction of nucleus to bonding / shared pair (of electrons) due to <i>any one of</i>:</p> <ul style="list-style-type: none"> more / high shielding (of electrons from inner shells) more / 5 electron shells in iodine C-I greater distance from nuclei (to bonding pair of electrons) | 2 |
| (c)(i) |  <p>M1 dipole C^{δ+}-Br^{δ-} AND curly arrow from C-Br bond to Br^{δ-}</p> <p>M2 lone pair on C of ⁻CN AND curly arrow from lone pair to C (on C-1 of bromopropane)</p> | 2 |

Q7.

| | | |
|----------|---|---|
| (e)(i) | <p>M1 correct dipole in on $\delta^+C-Br^{\delta-}$ (in 'BuBr) AND curly arrow from C—Br to Br</p> <p>M2 curly arrow from electron pair on O of $:OH^-$ / $:OH^-$ to C^+ in carbocation</p> <p>M3 'BuOH product</p> | 1 |
| (e)(ii) | The C—Br carbon is bonded to three other carbon atoms | 1 |
| (e)(iii) | | 1 |

Q8.

| | | |
|---------|---|---|
| (e)(i) | elimination | 1 |
| (e)(ii) | M1 NaOH / KOH | 1 |
| | M2 ethanolic solution / ethanol / alcohol + heat | 1 |

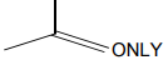
Q9.

| | | |
|---------|---|---|
| (c)(i) | <p>M1 correct dipole on haloalkane AND arrow from bond to Br or just beyond</p> <p>M2 correct intermediate</p> <p>M3 arrow from lone pair on <u>O</u> of OH^- / $-OH$ to central C of their intermediate shown in M2 OR arrow from lone pair on <u>O</u> of OH^- / $-OH$ to central C of 2-bromo 2-methylpropane if S_N2 mechanism shown</p> | 1 |
| (c)(ii) | nucleophilic substitution | 1 |
| (d) | M1 more time (because the rate is lower) | 1 |
| | M2 C—Cl (bond) is stronger (than C—Br) | 1 |

Q10.

| | | |
|----------|--|----------|
| (b)(iii) | <i>reagent</i> M1 NaCN or KCN | 1 |
| | <i>conditions</i> M2 ethanolic AND heat (under reflux) | 1 |

Q11.

| | | |
|----------|--|----------|
| (b)(i) | CH ₃ CH ₂ CH ₂ CH ₂ CH ₂ Br | 1 |
| (b)(i) | M1 primary/ 1° (carbo)cation formed is not very stable M2 EITHER (as) only one alkyl group exerting an inductive effect OR only one alkyl group so the charge is (more) localised on the C+ | 2 |
| (c)(i) | elimination | 1 |
| (c)(ii) | C ₄ H ₉ Cl + NaOH → C ₄ H ₈ + NaCl + H ₂ O | 1 |
| (c)(iii) |  ONLY | 1 |
| (c)(iv) | M1 2-chloro(-2-)methylpropane M2 1-chloro(-2-)methylpropane ALLOW in any order | 2 |

Q12.

| | | |
|----------|---|----------|
| (b)(i) | Human activity creates / additional / more/increase / thicker layer in greenhouse gas(es) / CHClF ₂ OR Human activity has an impact on climate change / temperature at earth's surface / temperature of sea | 1 |
| (b)(ii) | M1 traps (more)heat M2 (in the atmosphere leading to) greater global warming or wtte | 2 |
| (b)(iii) | ozone depletion / thinning | 1 |