M1.B

M2. (a) X contains $>\mathrm{C}=\mathrm{O}$ (1)
if $X$ and $Y$ reversed lose this mark but allow remaining max 6/7
$\therefore \mathrm{X}$ is $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}$ (1)
$\therefore \mathrm{Y}$ is $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{OH}$ (1)


(1)

Conc $\mathrm{H}_{2} \mathrm{SO}_{4}$ : catalyst (1)
(b)

(1)

B
 in any order
(c) $-\mathrm{O}_{\mathrm{C}}^{\mathrm{C}} \mathrm{H}_{2}-\quad 3.1-3.9(1)$

2.1-2.6(1)
a: quartet (1) $\quad 3$ adjacent $\mathrm{H}(1)$
b: triplet (1) $\overparen{2}$ adjacent $\mathrm{H}(\mathbf{1})$
(d) $3269 \mathrm{~cm}^{-1} \therefore \mathrm{OH} \curvearrowleft$ alcohol (1)

$\therefore \underline{\mathrm{G}}$ is
 (1)

## Notes

(a) first mark for $\mathrm{C}=\mathrm{O}$ stated or shown in X

Ignore wrong names
$\mathrm{Y} \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{OH}$
allow $\mathrm{C}_{3} \mathrm{H}_{7}$ in $\mathbf{A}$ if $\mathbf{Y}$ correct or vice versa
Allow (1) for $\mathbf{A}$ if correct conseq to wrong $\mathbf{X}$ and $\mathbf{Y}$
other oxidising agents: acidified $\mathrm{KMnO}_{4}$; Tollens; Fehlings
other reducing agents: $\mathrm{LiAlH}_{4} ; \mathrm{Na} /$ ethanol; $\mathrm{Ni} / \mathrm{H}_{2} ; \mathrm{Zn}$ or Sn or $\mathrm{Fe} / \mathrm{HCl}$
(b) give (1) for carboxylic acid stated or COOH shown in each suggestion (1) for correct E any 2 out of 3 for $\mathbf{B}, \mathbf{C}$ or $\mathbf{D}$ allow $\mathrm{C}_{3} \mathrm{H}_{7}$ for either the $\mathbf{B}$ or $\mathbf{D}$ shown on the mark scheme i.e. a correct structure labelled $\mathbf{B}, \mathbf{C}$ or $\mathbf{D}$ or $\mathbf{E}$ will gain 2.
(c) protons a - quartet must be correct to score 3 adjacent $H$ mark. Same for b
(d) allow (1) for any OH (alcohol) shown correctly in any structure - ignore extra functional groups. Structure must be completely correct to gain second mark

## Organic points

(1) Curly arrows: must show movement of a pair of electrons, i.e. from bond to atom or from Ip to atom / space
e.g.


OR

(2) Structures
penalise sticks (i.e.



or $-\mathrm{NH}_{2} \quad \checkmark$


$\mathrm{H}_{2} \mathrm{~N}-$
etc

## Penalise once per paper

$$
\frac{}{} \quad \frac{\mathrm{CH}_{3}}{\text { allow }} \mathrm{CH}_{3}-\text { or }-\mathrm{CH}_{3} \text { or } \stackrel{\mathrm{I}}{ }{ }^{\circ} \text { or } \mathrm{CH}_{3}
$$

M3.D

M4. (a) $\mathrm{CH}_{3} \mathrm{COCl}+\mathrm{AlCl}_{3} \rightarrow \mathrm{CH}_{3} \stackrel{+}{\mathrm{C}} \mathrm{O}+\mathrm{AICl}^{-}$
(1)
equation (1)
penalise wrong alkyl group once at first error position of + on electrophile can be on O or C or outside [ ]
penalise wrong curly arrow in the equation or lone pair on $\mathrm{AlCl}_{3}$ else ignore

Electrophilic substitution
NOT F/C acylation

horseshoe must not extend beyond C2 to C6 but can be smaller

+ not too close to C1
M3 arrow into hexagon unless Kekule
allow M3 arrow independent of M2 structure
M1 arrow from within hexagon to C or to + on C
+ must be on C of RCO
(b) Nucleophilic addition

NOT reduction


M3
M2 not allowed independent, but can allow M1 for attack of H on C+ formed

1-phenylethan(-1-)ol or (1-hydroxyethyl)benzene
(c) dehydration or elimination
(conc) $\mathrm{H}_{2} \mathrm{SO}_{4}$ or (conc) $\mathrm{H}_{3} \mathrm{PO}_{4}$
allow dilute and $\mathrm{Al}_{2} \mathrm{O}_{3}$
Do not allow iron oxides

M5. $\quad \mathbf{X}$ is $\mathrm{CH}_{3} \mathrm{CN}$ or ethanenitrile or ethanonitrile or methyl cyanide or cyanomethane or ethyl nitrile or methanecarbonitrile

Not ethanitrile
but contradiciton of name and structure lose marks
1
$\mathbf{Y}$ is $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{NH}_{2}$ or ethylamine or aminoethane or ethanamine

Step 1: reagent KCN not $\mathrm{HCN} / \mathrm{HCl}$ condition (aq)/alcohol - only allow condition if reagent correct or incomplete

| Step 2: reagent |  |  |  |  |  |
| ---: | :--- | :---: | :--- | :--- | :--- |
| condition | $\mathrm{H}_{2}$ | $\mathrm{NiAlH}_{4} / \mathrm{Pt} / \mathrm{Pd}$ | Na <br> ether | $\mathrm{Na} / \mathrm{Fe} / \mathrm{Sn}$ <br> ethanol | $\mathrm{Not} \mathrm{NaBH}_{4}$ |
| HCl |  |  |  |  |  |

$\mathbf{Z}$ is an amine or aminoalkane or named amine even if incorrect name for $\mathbf{Z}$ secondary (only award if amine correct)

$(\mathrm{Br})+$ can be on N or outside brackets as shown
nucleophilic substitution

M6.
(a) (i)

| Reagent | Tollens | Fehlings or Benedicts | $\mathrm{K}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{H}^{+}$ <br> or acidified | $\mathrm{KMnO}_{4} / \mathrm{H}^{+}$ | $\mathrm{l}_{2} / \mathrm{NaOH}$ |
| :--- | :--- | :--- | :--- | :--- | :--- |
| Propanal | silver <br> (mirror) | red ppt or goes red <br> (not red solution) | goes green | goes colourless | No <br> reaction |
| Propanone | no <br> reaction | no reaction | no reaction | no reaction | Yellow <br> (ppt) |

(penalise incomplete reagent e.g. $\mathrm{K}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7}$ or $\mathrm{Cr}_{2} \mathrm{O}_{7}^{2-} / \mathrm{H}^{+}$then mark on)
(ii) propanal 3 peaks ignore splitting even if wrong propanone 1 peak
(b) X is $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}$ or propanoic acid if both name and formula given, both must be correct, but
$\mathbf{Y}$ is $\mathrm{CH}_{3} \mathrm{CH}(\mathrm{OH}) \mathrm{CH}_{3}$ or propan-2-ol allow propanol with correct formula

Mark the type of reaction and reagent/condition independently. The reagent must be correct or close to score condition

Step 1 Oxidation
$\mathrm{K}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{H}^{+}$or other oxidation methods as above allow $\mathrm{Cr}_{2} \mathrm{O}_{7}^{2-} \mathrm{H}^{+}$if penalised above (ecf) reflux (not Tollens/Fehlings) or heat or warm

Step 2

| eduction or nucleophilic <br> addition | reduction or <br> nucleophilic addition | reduction or <br> hydrogenation |
| :--- | :--- | :--- |
| $\mathrm{NaBH}_{4}$ | LiAlH $_{4}$ | $\mathrm{H}_{2}$ |
| in (m)ethanol or water or <br> ether | ether or dry | $\mathrm{Ni} / \mathrm{Pt}$ etc |

1

1
$\qquad$

$\qquad$

