| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 ( a ) ( i )}$ | Concentrated nitric acid AND <br> concentrated sulfuric acid <br> ALLOW <br> 'concentrated nitric and sulfuric <br> acids' <br> Concentrated $\mathrm{HNO}_{3}$ and concentrated <br> $\mathrm{H}_{2} \mathrm{SO}_{4}$ | Extra reagents |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 ( a ) ( i i )}$ | To prevent multiple substitutions/ to <br> stop di- or trinitrobenzene forming <br> ALLOW <br> To stop further substitution (of $\mathrm{NO}_{2}$ )/ <br> further nitration <br> IGNORE <br> further reaction | Further addition <br> of nitro groups | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 ( a ) ( \text { iii) }}$ | Tin/ Sn AND concentrated HCl/ <br> concentrated hydrochloric acid | Dilute HCl | 1 |
|  | ALLOW Iron/Fe or Zn/Zinc for tin <br> Conc for concentrated |  | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 ( b ) ( i )}$ | $\mathrm{C}_{6} \mathrm{H}_{5} \mathrm{NH}_{3}{ }^{+} \mathrm{Cl}^{-}$ |  | 1 |
|  | $\mathrm{ALLOW}^{\mathrm{C}_{6} \mathrm{H}_{5} \mathrm{NH}_{3} \mathrm{Cl}}$ |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 1(b)(ii) |   <br> H <br> ALLOW <br> $\mathrm{C}_{6} \mathrm{H}_{5}$ for benzene <br> Undisplayed $\mathrm{CH}_{3}$ | Skeletal formula <br> Structural formula | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 ( b ) ( i i i ) i}$ | (transition metal) complex ion <br> ALLOW <br> Transition metal complex / copper <br> complex <br> IGNORE (1) <br> Formulae of ions |  | 2 |
|  | F (azo) dye / azo compound / diazo <br> compound <br> ALLOW <br> diazonium compound <br> molecule for compound |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 b ( i v ) ~}$ | Benzenediazonium chloride <br> ALLow <br> Phenyldiazonium chloride | Benzadiazonium <br> chloride <br> Diazonium salt | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 b ( v )}$ | $\mathrm{HCl}+\mathrm{NaNO}_{2}$ <br> OR <br> Hydrochloric acid + Sodium nitrite / <br> nitrate(III) <br> OR <br> alternative cation to $\mathrm{Na}^{+}$ | $\mathrm{HCl}+\mathrm{HNO}_{2}$ | 1 |
| IGNORE <br> $\mathrm{HNO}_{2}$ <br> Concentration of HCl |  |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 1b(vi) |  <br> ALLOW <br> any substitution positions $\mathrm{C}_{6} \mathrm{H}_{3}\left(\mathrm{CH}_{3}\right)_{2} \mathrm{NH}_{2}$ <br> H- ${ }_{6} \mathrm{H}_{2}\left(\mathrm{CH}_{3}\right)_{2} \mathrm{NH}_{2}$ <br> Kekule structure | $\mathrm{C}_{6} \mathrm{H}_{\mathbf{2}}\left(\mathrm{CH}_{3}\right)_{2} \mathrm{NH}_{2}$ | 1 |

Total for Question = 10 marks

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 ~ ( a ) ( i ) ~}$ | Overall yield higher <br> OR <br> Reduces use of solvents (ALLOW <br> chemicals / reactants) <br> OR <br> Less loss of chemicals <br> OR <br> Less waste products <br> IGNORE <br> References to Energy / fuel / $\mathrm{CO}_{2}$ <br> References to atom economy <br> More efficient conversion <br> Fewer side products | $\mathbf{1}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2}$ | Lowers (operating) temperature / <br> energy (requirements) <br> OR <br> Less fuel needed <br> (a)(ii) | IGNORE <br> References to catalyst properties such <br> as 'lowers E ${ }_{\text {a ', 'can be re-used' }}$ <br> Atom economy | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 ~ ( b ) ( i ) ~}$ | $\mathrm{CH}_{3} \mathrm{COCl}+\mathrm{AlCl}_{3} \rightarrow \mathrm{CH}_{3} \mathrm{CO}^{+}+\mathrm{AlCl}_{4}{ }^{-}$Structural formulae not required <br> Positive charge may be anywhere on <br> the electrophile. <br> IGNORE <br> Curly arrows even if incorrect | $\mathbf{1}$ |  |


| Question |
| :--- | :--- | :--- | :--- |
| Number | Acceptable Answers


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2}$ (b)(iii) | No HCl formed (as a by-product) <br> OR <br> Ethanoic acid easier to recover <br> ALLOW <br> Reverse arguments <br> IGNORE <br> Chlorine containing product <br> References to ozone layer, acid rain, <br> global warming <br> Atom economy | Chlorine | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 ( c ) ( i )}$ | Catalyst (more) easily recovered / <br> separated OR can be filtered |  | $\mathbf{1}$ |
| OR <br> Facilitates the use of flow (rather than <br> batch) systems <br> IGNORE <br> references to properties of catalysts |  |  |  |



| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 2 <br> (c) (iii) | $\mathrm{C}=\mathrm{O} /$ carbonyl group (only) in carboxylic acid / ibuprofen <br> Absorption / peak <br> at 1725-1700 ( $\mathrm{cm}^{-1}$ ) <br> If no other mark has been awarded, then ALLOW (for 1 mark) <br> OH in both but in alcohol 3750-3200 $\left(\mathrm{cm}^{-1}\right)$ but in | ketone $\begin{align*} & 1700-1680  \tag{1}\\ & \left(\mathrm{~cm}^{-1}\right) \end{align*}$ <br> Single values rather than ranges | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 ~ ( d ) ( i )}$ | (A chiral molecule is) non-superimposable on its <br> mirror image. |  | $\mathbf{1}$ |
|  | ALLOW <br> Asymmetric (tetrahedral) carbon atom / has a <br> carbon atom bonded to four different groups / <br> atoms | molecules <br> / species <br> (for <br> groups) | IGNORE <br> Has two enantiomers <br> Functional (as in functional groups) <br> Reference to rotation of plane polarized light |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2}$ |  |  |  |
| $\mathbf{( d ) ( i i )}$ |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2}$ | (d) racemic mixture is) an equimolar <br> (dixture of the two enantiomers / (optical) <br> isomers <br> ALLOW (for equimolar mixture) <br> equal amounts / concentrations / volumes / <br> proportions | Just 'no effect on <br> plane polarised <br> light' | $\mathbf{1}$ |
| OR <br> $50: 50$ mixture |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2}$ | Any two of |  | $\mathbf{2}$ |
| (d)(iv) | 1. All the ibuprofen is useful (rather than half) <br> 2.No need for separation of isomers / <br> enantiomers <br> 3. No need for a more complex synthesis <br> forming just one enantiomer <br> 4. Sometimes one enantiomer has negative <br> effects | 5. Smaller dosage may be used |  |
| ALLOW (For point 4 above) <br> Dose / inactive isomer is less likely to be <br> harmful | IGNORE <br> Reference to cost / yield / atom economy / side <br> effects |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{3}$ <br> (a)(i) | $\mathrm{HNO}_{3}+2 \mathrm{H}_{2} \mathrm{SO}_{4} \rightarrow \mathrm{H}_{3} \mathrm{O}^{+}+2 \mathrm{HSO}_{4}{ }^{-}+\mathrm{NO}_{2}{ }^{+}$ |  |  |
| OR |  | 2 |  |
|  | $\mathrm{HNO}_{3}+\mathrm{H}_{2} \mathrm{SO}_{4} \rightarrow \mathrm{H}_{2} \mathrm{O}+\mathrm{HSO}_{4}{ }^{-}+\mathrm{NO}_{2}{ }^{+}$ <br> OR <br> 2-step version of these involving $\mathrm{H}_{2} \mathrm{NO}_{3}{ }^{+}$ <br> Correct electrophile (1) <br> correct equation(s) (1) |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{gathered} 3 \\ \text { (a) }(\mathrm{ii}) \end{gathered}$ | OR $\mathrm{NO}_{2}{ }^{+}$as electrophile <br> TE on incorrect electrophile in (a)(i) <br> Curly arrow from on or within the circle to positively charged nitrogen <br> ALLOW <br> Curly arrow from anywhere within the hexagon <br> Arrow to any part of the electrophile including to the + charge <br> Intermediate structure including charge with horseshoe covering at least 3 carbon atoms, and <br> facing the tetrahedral carbon and <br> with some part of the positive charge within the horseshoe <br> Curly arrow from $\mathrm{C}-\mathrm{H}$ bond to anywhere in the benzene ring reforming delocalized structure <br> Correct Kekulé structures score full marks <br> Ignore any involvement of anion in the final step |  | 3 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{3}$ | Benzene ring in phenol has higher <br> electron density <br> (a)(iii) <br> O / OH donates electron density to <br> the (benzene) ring (1) <br> Because lone pair of electrons on <br> (phenol) oxygen is donated to / <br> overlaps with / interacts with ( $\square$ <br> electrons of benzene) ring (1) |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{3}$ | Substitution may also occur at the 2 / <br> (a)(iv) <br> 6 ring positions / ortho position |  | 1 |
|  | ALLOW 'other' / 3 / 5 / meta ring <br> positions / isomers |  |  |
|  | ALLOW further substitution occurs |  |  |
| IGNORE |  |  |  |
| By-products formed |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{3}$ | Tin /Sn \& (conc.) hydrochloric acid / | $\mathrm{LiAlH}_{4} / \mathrm{NaBH}_{4}$ | 1 |
| $\mathbf{( a ) ( v )}$ | Cl(aq) <br> ALLOW <br> Iron/ Fe for tin <br> ALLOW HCl for HCl(aq) |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{3}$ | Yield $=(100 \times 0.25 \times 0.74 \times 0.85)=$ <br> $(\mathbf{a})(\mathbf{v i})$ | 16.0 and other <br> rounding errors | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{3 ( b ) ( i )}$ | Insoluble impurities are removed by <br> the hot filtration <br> Soluble impurities are removed by <br> the cold filtration |  | 2 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{gather*} 3  \tag{1}\\ \text { (b) }(\mathrm{ii}) \end{gather*}$ | 5으 and 95응 <br> Because the lowest proportion (ALLOW 'amount') of paracetamol remains in solution (at the end) <br> IGNORE <br> Just 'greatest difference in temperature' |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{3}$ (b)(iii) | Measure melting temperature | Boiling <br> temperature | 1 |
|  | ALLOW <br> TLC (with UV light) <br> Ignore <br> Must melt over range of 2으 <br> Data = data book value | HPLC |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{3 ~ c ( i ) ~}$ | Peak at $\mathrm{m} / \mathrm{e}=151$ clearly labelled $\mathbf{M}$ <br>  <br>  <br> ALLOW <br> Alternative labels | 1 |  |


| Question Number | Acceptable Answers |  | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 3 c (ii) | $43=\left[\mathrm{CH}_{3}-\mathrm{C}{ }_{\mathrm{O}}\right]^{+}$ <br> ALLOW CONH ${ }^{+}$ <br> Ignore position of charges | $\begin{gathered} \mathrm{OR} \\ \mathrm{CH}_{3} \mathrm{CO}^{+} \\ \mathrm{C}_{2} \mathrm{H}_{3} \mathrm{O}^{+} \end{gathered}$ | $\mathrm{C}_{3} \mathrm{H}_{7}^{+}$ <br> uncharged species | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{3}$ (d) | Limit number of tablets sold <br> OR <br> Give (oral) advice at the point of sale <br> OR <br> Use packs with tablets individually <br> wrapped <br> ALLOW <br> Reduce the (tablet) dose | Only sell on <br> prescription / <br> doctor's advice <br> Label packet | 1 |

