

AQA Qualifications

## A-level **Chemistry**

Paper 2 (7405/2): Organic and Physical Chemistry Mark scheme

7405 Specimen paper

Version 0.5

| Question | Marking guidance                                                                                                                           | Mark | AO     | Comments                                               |
|----------|--------------------------------------------------------------------------------------------------------------------------------------------|------|--------|--------------------------------------------------------|
| 01.1     | Consider experiments 1 and 2: [B constant]                                                                                                 |      |        |                                                        |
|          | [A] increases × 3: rate increases by 3 <sup>2</sup> therefore 2nd order with respect to A                                                  | 1    | AO3 1a |                                                        |
|          | Consider experiments 2 and 3:                                                                                                              |      |        |                                                        |
|          | [A] increases × 2: rate should increase × 2 <sup>2</sup> but only increases × 2                                                            |      |        |                                                        |
|          | Therefore, halving [B] halves rate and so 1st order with respect to B                                                                      | 1    | AO3 1a |                                                        |
|          | Rate equation: $rate = k[A]^2[B]$                                                                                                          | 1    | AO3 1b |                                                        |
| 01.2     | rate = $k [C]^2 [D]$ therefore $k = \text{rate} / [C]^2 [D]$                                                                               | 1    | AO2h   |                                                        |
|          | $k = \frac{7.2 \times 10^{-4}}{(1.9 \times 10^{-2})^2 \times (3.5 \times 10^{-2})} = 57.0$                                                 | 1    | AO2h   | Allow consequential marking on incorrect transcription |
|          | $mol^{-2} dm^{+6} s^{-1}$                                                                                                                  | 1    | AO2h   | Any order                                              |
| 01.3     | rate = $57.0 \times (3.6 \times 10^{-2})^2 \times 5.4 \times 10^{-2} = 3.99 \times 10^{-3} \text{ (mol dm}^{-3} \text{ s}^{-1})$ <b>OR</b> | 1    | AO2h   |                                                        |
|          | Their $k \times (3.6 \times 10^{-2})^2 \times 5.4 \times 10^{-2}$                                                                          |      |        |                                                        |

| 01.4 | Reaction occurs when molecules have $E \ge E_a$ Doubling T causes many more molecules to have this $E$ Whereas doubling [E] only doubles the number with this $E$ | 1<br>1<br>1 | AO1a<br>AO1a<br>AO1a |                                                                                              |
|------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------|-------------|----------------------|----------------------------------------------------------------------------------------------|
| 01.5 | $E_a = RT(\ln A - \ln k)/1000$                                                                                                                                    | 1           | AO1b                 | Mark is for rearrangement of equation and factor of 1000 used correctly to convert J into kJ |
|      | $E_{\rm a} = 8.31 \times 300 (23.97 - (-5.03))/1000 = 72.3 (kJ mol^{-1})$                                                                                         | 1           | AO1b                 |                                                                                              |

| Question | Marking guidance                                                                                                                                                                                         | Mark   | AO     | Comments                                                         |
|----------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|--------|--------|------------------------------------------------------------------|
| 02.1     | Marking guidance  Gradient drawn on graph  0.020 0.018 0.016 0.014 0.012  [Butadiene] / mol dm <sup>-3</sup> 0.010 0.008 0.006 0.004 0.002 0.000 0 1000 2000 3000 4000 5000 6000 7000 8000 9000 Time / s | Mark 1 | AO3 1a | Line must touch the curve at 0.012 but must not cross the curve. |
|          | Time/s                                                                                                                                                                                                   |        |        |                                                                  |

| 02.2 |                                                                                                                                                                                                                          |   |                  | Extended response                                |
|------|--------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|---|------------------|--------------------------------------------------|
| 02.2 | Stage 1: Rate of reaction when concentration = 0.0120 mol dm <sup>-3</sup> From the tangent                                                                                                                              |   |                  | Extended response                                |
|      | Change in [butadiene] = $-0.0160 - 0$ and change in time = $7800 - 0$<br>Gradient = $-(0.0160 - 0)/(7800 - 0) = -2.05 \times 10^{-6}$                                                                                    | 1 | AO3 1a           |                                                  |
|      | Rate = $2.05 \times 10^{-6}$ (mol dm <sup>-3</sup> s <sup>-1</sup> )                                                                                                                                                     | 1 | AO3 1a           |                                                  |
|      | Stage 2: Comparison of rates and concentrations Initial rate/rate at $0.0120 = (4.57 \times 10^{-6})/(2.05 \times 10^{-6}) = 2.23$ Inital concentration/concentration at point where tangent drawn = $0.018/0.012 = 1.5$ | 1 | AO3 1a<br>AO3 1a | Marking points in stage 2 can be in either order |
|      | Stage 3: Deduction of order  If order is 2, rate should increase by factor of $(1.5)^2 = 2.25$ this is approximately equal to 2.23 therefore order is 2nd with respect to butadiene                                      | 1 | AO3 1b           |                                                  |

| Question | Marking guidance                                             | Mark | AO   | Comments |
|----------|--------------------------------------------------------------|------|------|----------|
| 03.1     | 2,2,4-trimethylpentane                                       | 1    | AO1a |          |
| 03.2     | 5                                                            | 1    | AO2b |          |
| 03.3     | $C_{20}H_{42} \longrightarrow C_8H_{18} + 2C_3H_6 + 3C_2H_4$ | 1    | AO2b |          |
| 03.4     | Mainly alkenes formed                                        | 1    | AO1b |          |
| 03.5     | 4 (monochloro isomers)                                       | 1    | AO2b |          |
|          | $CH_3 H CH_3$<br>                                            | 1    | AO2a |          |
| 03.6     | CI                                                           | 1    | AO2a |          |

| 03.7 | $C_8H_{17}^{35}CI = 96.0 + 17.0 + 35.0 = 148.0$<br>and $C_8H_{17}^{37}CI = 96.0 + 17.0 + 37.0 = 150.0$             | 1 | AO1b | Both required |
|------|--------------------------------------------------------------------------------------------------------------------|---|------|---------------|
|      | $M_{\rm r}$ of this C <sub>8</sub> H <sub>17</sub> Cl $(1.5 \times 148.0) + (1.0 \times 150.0) = 148.8$<br>2.5 2.5 | 1 | AO1b |               |
| 03.8 | <u>24.6</u> <u>2.56</u> <u>72.8</u> = 2.05 : 2.56 : 2.05<br>12 1 35.5                                              |   |      |               |
|      | Simplest ratio = $\frac{2.05}{2.05}$ : $\frac{2.56}{2.05}$ : $\frac{2.05}{2.05}$                                   |   |      |               |
|      | = 1 : 1.25 : 1                                                                                                     | 1 | AO2b |               |
|      | Whole number ratio ( $\times$ 4) = 4 : 5 : 4                                                                       | 1 | AO2b |               |
|      | $MF = C_8 H_{10} CI_8$                                                                                             | 1 | AO2b |               |

| Question | Marking guidance                                             | Mark | AO   | Comments                                                   |
|----------|--------------------------------------------------------------|------|------|------------------------------------------------------------|
| 04.1     | 3-methylbutan-2-ol                                           | 1    | AO1a |                                                            |
| 04.2     | CH <sub>3</sub> H <sub>3</sub> C—C—C—CH <sub>3</sub> H O     | 1    | AO2g | Allow (CH <sub>3</sub> ) <sub>2</sub> CHCOCH <sub>3</sub>  |
| 04.3     | Elimination                                                  | 1    | AO1a |                                                            |
| 04.4     | CH <sub>3</sub>   H <sub>3</sub> C—C=C-CH <sub>3</sub>   H   | 1    | AO2g | Allow (CH <sub>3</sub> ) <sub>2</sub> C=CHCH <sub>3</sub>  |
|          | CH <sub>3</sub>   H <sub>3</sub> C—C—C—CH <sub>2</sub>   H H | 1    | AO2g | Allow (CH <sub>3</sub> ) <sub>2</sub> CHCH=CH <sub>2</sub> |

| 04.5 | Position                                             | 1 | AO1a   |                                                                            |
|------|------------------------------------------------------|---|--------|----------------------------------------------------------------------------|
| 04.6 | СВА                                                  | 1 | AO3 1b |                                                                            |
| 04.7 | H <sub>3</sub> C——CH <sub>2</sub> CH <sub>3</sub> OH | 1 | AO2g   | Allow (CH <sub>3</sub> ) <sub>2</sub> C(OH)CH <sub>2</sub> CH <sub>3</sub> |
| 04.8 | $H_3C$ — $C$ — $CH_2OH$ $CH_3$                       | 1 | AO2e   | Allow (CH <sub>3</sub> ) <sub>3</sub> CCH <sub>2</sub> OH                  |

| Question | Marking guidance                                                                                                                                                                                                                                   | Mark             | AO                           | Comments |
|----------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|------------------|------------------------------|----------|
| 05.1     | Secondary                                                                                                                                                                                                                                          | 1                | AO1a                         |          |
| 05.2     | Nitrogen and oxygen are very electronegative Therefore, C=O and N-H are polar Which results in the formation of a hydrogen bond between O and H In which a lone pair of electrons on an oxygen atom is strongly attracted to the $\delta\text{+H}$ | 1<br>1<br>1<br>1 | AO1a<br>AO1a<br>AO1a<br>AO1a |          |

| Question | Marking guidance                                      | Mark | AO   | Comments                                                  |
|----------|-------------------------------------------------------|------|------|-----------------------------------------------------------|
| 06.1     | $\begin{array}{c ccccccccccccccccccccccccccccccccccc$ | 1    | AO2a |                                                           |
| 06.2     | +                                                     | 1    | AO2a |                                                           |
| 06.3     | $\begin{array}{c ccccccccccccccccccccccccccccccccccc$ | 1    | AO2a | Allow $(CH_3)_3$ $\stackrel{+}{N}$ $-CH_2$ $-COOH$ $(Br)$ |
| 06.4     | 2-amino-3-hydroxybutanoic acid                        | 1    | AO2a |                                                           |

| 06.5 | +<br>ŅH <sub>3</sub>                | 1 | AO2a |  |
|------|-------------------------------------|---|------|--|
|      | <br>(CH <sub>2</sub> ) <sub>4</sub> |   |      |  |
|      | H <sub>3</sub> N—C—COOH             |   |      |  |
|      | н                                   |   |      |  |

| Question | Marking guidance                                                                  | Mark | AO   | Comments |
|----------|-----------------------------------------------------------------------------------|------|------|----------|
| 07.1     | H CH <sub>3</sub>                                                                 | 1    | AO1a |          |
|          | Ċ──Ċ<br>I I<br>CH <sub>3</sub> CI                                                 |      |      |          |
|          | Addition                                                                          | 1    | AO1a |          |
| 07.2     | Ų Ų                                                                               | 1    | AO2e |          |
|          | H H<br>  HO-C-C-OH<br>  CH <sub>3</sub> CH <sub>3</sub>                           |      |      |          |
|          | O CH <sub>3</sub> H O O CH <sub>3</sub> H O II I II | 1    | AO2e |          |
| 07.3     | <b>Q</b> is biodegradable                                                         | 1    | AO2g |          |
|          | Polar C=O group or $\delta$ + C in <b>Q</b> (but not in <b>P</b> )                | 1    | AO2c |          |
|          | Therefore, can be attacked by nucleophiles (leading to breakdown)                 | 1    | AO2c |          |

| Question | Marking guidance                                                                                                                                                                     | Mark             | AO                             | Comments                                                                                                                                                                                                                |
|----------|--------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|------------------|--------------------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| 08.1     | 2-deoxyribose                                                                                                                                                                        | 1                | AO1a                           |                                                                                                                                                                                                                         |
| 08.2     | Base A Top N–H forms hydrogen bonds to lone pair on O of guanine The lone pair of electrons on N bonds to H–N of guanine A lone pair of electrons on O bonds to lower H–N of guanine | 1<br>1<br>1<br>1 | AO3 1b<br>AO2a<br>AO2a<br>AO2a | If Base B stated, allow 1 mark only for response including hydrogen bonding  Allow all 4 marks for a correct diagram showing the hydrogen bonding  Students could also answer this question using labels on the diagram |
| 08.3     | Allow either of the nitrogen atoms with a lone pair NOT involved in bonding to cytosine                                                                                              | 1                | AO2a                           |                                                                                                                                                                                                                         |
| 08.4     | Use in very small amounts / target the application to the tumour                                                                                                                     | 1                | AO2e                           |                                                                                                                                                                                                                         |

| Question | Marking guidance                                     | Mark | AO   | Comments                                                                                                                                                                                                                                                                                    |
|----------|------------------------------------------------------|------|------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| 09.1     | (nucleophilic) addition-elimination                  | 1    | AO1a | Not electrophilic addition-elimination  Allow C <sub>6</sub> H <sub>5</sub> or benzene ring                                                                                                                                                                                                 |
|          | $\begin{array}{cccccccccccccccccccccccccccccccccccc$ | 4    | AO2a | Allow attack by :NH <sub>2</sub> C <sub>6</sub> H <sub>5</sub> M2 not allowed independent of M1, but allow M1 for correct attack on C+  M3 for correct structure with charges but lone pair on O is part of M4  M4 (for three arrows and lone pair) can be shown in more than one structure |

| 09.2 | The minimum quantity of hot water was used:                                    |   |        |                                                |
|------|--------------------------------------------------------------------------------|---|--------|------------------------------------------------|
|      | To ensure the hot solution would be saturated / crystals would form on cooling | 1 | AO1b   |                                                |
|      | The flask was left to cool before crystals were filtered off:                  |   |        |                                                |
|      | Yield lower if warm / solubility higher if warm                                | 1 | AO1b   |                                                |
|      | The crystals were compressed in the funnel:                                    |   |        |                                                |
|      | Air passes through the sample not just round it                                | 1 | AO1b   | Allow better drying but not water squeezed out |
|      | A little cold water was poured through the crystals:                           |   |        |                                                |
|      | To wash away soluble impurities                                                | 1 | AO1b   |                                                |
| 09.3 | Water                                                                          | 1 | AO3 1b | Do not allow unreacted reagents                |
|      | Press the sample of crystals between filter papers                             | 1 | AO3 2b | Allow give the sample time to dry in air       |
| 09.4 | $M_{\rm r}$ product = 135.0                                                    | 1 | AO2h   |                                                |
|      | Expected mass = $5.05 \times \frac{135.0}{93.0} = 7.33 \text{ g}$              | 1 | AO2h   |                                                |
|      | Percentage yield = $\frac{4.82}{7.33} \times 100 = 65.75 = 65.8(\%)$           | 1 | AO1b   | Answer must be given to this precision         |

| 09.5 | NHCOCH <sub>3</sub> $+ NO_2^+ \longrightarrow NHCOCH_3$ $+ NO_2^+ \longrightarrow NO_2$ OR $C_6H_5NHCOCH_3 + NO_2^+ \rightarrow C_6H_4(NHCOCH_3)NO_2 + H^+$ | 1 | AO2c   |                                         |
|------|-------------------------------------------------------------------------------------------------------------------------------------------------------------|---|--------|-----------------------------------------|
| 09.6 | Electrophilic substitution                                                                                                                                  | 1 | AO1a   |                                         |
| 09.7 | Hydrolysis                                                                                                                                                  | 1 | AO3 1a |                                         |
| 09.8 | Sn/HCI                                                                                                                                                      | 1 | AO1b   | Ignore acid concentration; allow Fe/HCI |

| Question |          | Marking guidance                                                                                | Mark | AO     | Comments                                                                                           |
|----------|----------|-------------------------------------------------------------------------------------------------|------|--------|----------------------------------------------------------------------------------------------------|
| 10       | IR<br>M1 | Absorption at 3360 cm <sup>-1</sup> shows OH alcohol present                                    | 1    | AO3 1a | Extended response  Deduction of correct structure without explanation                              |
|          | NMR      |                                                                                                 | ·    |        | scores maximum of 4 marks as this does not show a clear, coherent line of reasoning.               |
|          | M2       | There are 4 peaks which indicates 4 different environments of hydrogen                          | 1    | AO3 1a | Maximum of 6 marks if no structure given OR if coherent logic not displayed in the explanations of |
|          | M3       | The integration ratio = 1.6 : 0.4 : 1.2 : 2.4                                                   | 1    | AO3 1a | how two of OH, CH <sub>3</sub> and CH <sub>2</sub> CH <sub>3</sub> are identified.                 |
|          |          | The simplest whole number ratio is 4 : 1 : 3 : 6                                                |      |        |                                                                                                    |
|          | M4       | The singlet (integ 1) must be caused by H in OH alcohol                                         | 1    | AO3 1a |                                                                                                    |
|          | M5       | The singlet (integ 3) must be due to a CH <sub>3</sub> group with no adjacent H                 | 1    | AO3 1b |                                                                                                    |
|          | M6       | Quartet + triplet suggest CH <sub>2</sub> CH <sub>3</sub> group                                 | 1    | AO3 1b |                                                                                                    |
|          | M7       | Integration 4 and integration 6 indicates two equivalent CH <sub>2</sub> CH <sub>3</sub> groups | 1    | AO3 1b |                                                                                                    |
|          |          | ÇH₂CH₃                                                                                          |      |        |                                                                                                    |
|          | M8       | H <sub>3</sub> C——С—ОН                                                                          | 1    | AO3 1b |                                                                                                    |
|          |          | CH <sub>2</sub> CH <sub>3</sub>                                                                 |      |        |                                                                                                    |

| Question | Marking guidance                                           |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      | Mark | AO                     | Comments                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   |
|----------|------------------------------------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|------|------------------------|--------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| 11.1     | $CH_3CH_2COCH_3 + 2[H] \longrightarrow CH_3CH_2CH(OH)CH_3$ |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      |      | AO1b                   |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
| 11.2     |                                                            | All stages are covered and the explanation of each stage is generally correct and virtually complete.  Answer is communicated coherently and shows a logical progression from stage 1 to stage 2 then stage 3.  All stages are covered but the explanation of each stage may be incomplete or may contain inaccuracies OR two stages are covered and the explanations are generally correct and virtually complete.  Answer is mainly coherent and shows progression from stage 1 to stage 3.  Two stages are covered but the explanation of each stage may be incomplete or may contain inaccuracies, OR only one stage is covered but the explanation is generally correct and virtually complete.  Answer includes isolated statements but these are not presented in a logical order or show confused reasoning.  Insufficient correct chemistry to gain a mark. | 6    | 1<br>AO1a<br>5<br>AO2a | Indicative Chemistry content  Stage 1: Formation of product  Nucleophilic attack Planar carbonyl group Hattacks from either side (stated or drawn)  Stage 2: Nature of product Product of step 1 shown This exists in two chiral forms (stated or drawn) Equal amounts of each enantiomer/racemic mixture formed  Stage 3: Optical activity Optical isomers/enantiomers rotate the plane of polarised light equally in opposite directions With a racemic/equal mixture the effects cancel |

| Question | Marking guidance                                     | Mark | АО   | Comments                                                        |
|----------|------------------------------------------------------|------|------|-----------------------------------------------------------------|
| 12.1     | HBr OR HCI OR H <sub>2</sub> SO <sub>4</sub>         | 1    | AO1b | Allow HI or HY                                                  |
| 12.2     | Electrophilic addition                               | 1    | AO1a |                                                                 |
|          | $\begin{array}{cccccccccccccccccccccccccccccccccccc$ | 4    | AO2a | Allow consequential marking on acid in 12.1 and allow use of HY |
| 12.3     | The major product exists as a pair of enantiomers    | 1    | AO2a |                                                                 |
|          | The third isomer is 1-bromobutane (minor product)    | 1    | AO2a |                                                                 |
|          | Because it is obtained via primary carbocation       | 1    | AO2a |                                                                 |