| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (a) | (i) | $0 \checkmark,+2 \checkmark$ | 2 | NOT 2+ |
| 1 | (a) | (ii) | nitrogen(IV) oxide $\checkmark$ | 1 | ACCEPT nitrogen dioxide (or with no gap between words) or nitrogen(IV) dioxide (with or without gaps) <br> ALLOW nitrogen (IV) oxide ALLOW without brackets round 'IV' |
| 1 | (b) |  | brown (gas) (formed) $\checkmark$ | 1 | IGNORE 'colourless' as starting colour |
| 1 | (c) |  | advantage: nitrogen is fixed <br> OR it (enters the soil and) benefits plants (AW) $\checkmark$ <br> disadvantage: Acid rain OR an effect of acid rain eg buildings corroded / increasing acidity of soil / kills trees, increasing acidity of lakes (or rivers) / kills fish OR toxic <br> OR causes respiratory problems <br> OR (tropospheric) ozone formed <br> OR contributes to photochemical smog $\checkmark$ | 2 | idea of benefit to plants (eg: needed by plants / help plants grow / fertiliser / provides nutrients to soil, etc) needed for the 'or' version of the mark <br> IGNORE 'harmful'/'polluting' for last two points must have idea of giving rise to ozone or photochemical smog IGNORE 'greenhouse gas' |
| 1 | (d) | (i) | $0.2 / 24 \checkmark$ <br> OR $8.3 \times 10^{-3} \times 24=0.2$ AND comment that this is the fraction of oxygen in air <br> $0.033(2)$ or 0.033333 (to 2 or more sf) [or standard form] | 2 | Numbers 0.2 (or 1/5) and divided by 24 must be there OR 1/120 OR 2/240 <br> $8.3 \times 10^{-3}$ alone obviously does not score <br> ie do not accept '0.03’ <br> Answer alone scores the mark. Working alone does not score <br> ALLOW standard form eg $3.3 \times 10^{-2}$ ACCEPT recurring decimal <br> NOT rounding errors <br> watch out for $3.3 \times 10^{-3}$ (incorrect) |
| 1 | (d) | (ii) | $\left[\mathrm{NO}_{2}\right]^{2} /\left[\mathrm{N}_{2}\right]\left[\mathrm{O}_{2}\right]^{2} \checkmark$ | 1 | MUST have square brackets NOT ' p ' (even if with square brackets) ALLOW $\left[\mathrm{N}_{2}\right] \times\left[\mathrm{O}_{2}\right]^{2}$ or $\left[\mathrm{N}_{2}\right] \cdot\left[\mathrm{O}_{2}\right]^{2}$ NOT $\left[\mathrm{N}_{2}\right]+\left[\mathrm{O}_{2}\right]^{2}$ <br> ALLOW ' $(\mathrm{g})$ ' as state symbols. Others are CON |


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| 1 | (d) | (iii) | $\sqrt{ }\left(\left\{\right.\right.$ ans for $\left[\mathrm{N}_{2}\right]$ from d(i) $\left.) \times\left(8.3 \times 10^{-3}\right)^{2} \times 4 \times 10^{-19}\right) \checkmark$ OR $\sqrt{ }\left(\left[\mathrm{N}_{2}\right] \times\left[\mathrm{O}_{2}\right]^{2} \times K_{(c)}\right)$ OR part numbers part symbols evaluation (eg $\left[\mathrm{N}_{2}\right]=0.033$ gives $9.5(4) \times 10^{-13}$ ) $\checkmark$ ( 0.0332 gives $9.56 / 9.6 \times 10^{-13}$ ) <br> one sig fig $\left(e g 1 \times 10^{-12}\right) \checkmark$ | 3 | correct evaluation of correct expression (even if expression is not written down) scores 2 <br> No ecf from d(ii); no ecf for evaluating an incorrect expression, except that omission of the square root and then correctly evaluated (eg $\left.9(.1) \times 10^{-25}\right)$ or omission of square on $\mathrm{O}_{2}\left(\mathrm{eg} 1(.1) \times 10^{-11}\right)$ score one mark. <br> A spreadsheet is available for other answers from (d)(i) Mark sf separately, awarding the mark for any number to one sf. So proceed as follows: <br> - check for the three possible answers (with ecf if necessary using spreadsheet) One is worth two, others are worth one <br> - If no matching answer is present, look for square root expression and award 1 if it is correct <br> - Award sf mark if relevant |
| 1 | (d) | (iv) | $\mathrm{K}_{\mathrm{c}}$ larger $\checkmark$ <br> (Forward) reaction endothermic <br> OR right is endothermic direction ORA $\checkmark$ <br> Equilibrium position moves to right/towards products $\checkmark$ | 3 | Mark separately No ecf from second marking point incorrect IGNORE references to rates <br> must mention 'position' in connection with equilibrium |
| 1 | (d) | (v) | Equilibrium (position) moves to right/products <br> More molecules/moles on left-hand side/reactants ORA | 2 | Incorrect effects on $K_{\mathrm{c}}$ are CON of first mark (IGNORE 'no effect') <br> ALLOW 'it' for 'equilibrium position' <br> IGNORE 'favours the right-hand side' <br> ALLOW 'particles' instead of 'molecules' <br> IGNORE 'more reactants than products' (ie without mention of moles etc) <br> Mark separately |
| 1 | (e) | (i) | ```\(\mathrm{HNO}_{3} \rightarrow \mathrm{H}^{+}+\mathrm{NO}_{3}^{-}\) or \(\mathrm{HNO}_{3}+\mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{H}_{3} \mathrm{O}^{+}+\mathrm{NO}_{3}^{-}\) \(\checkmark\) for correct species \(\checkmark\) for correct species with arrow (not equm sign)``` | 2 | IGNORE state symbols <br> ALLOW $\mathrm{HNO}_{3}+\mathrm{aq} \rightarrow \mathrm{H}^{+}+\mathrm{NO}_{3}{ }^{-}$ <br> $\mathrm{HNO}_{3} \rightarrow \mathrm{H}_{3} \mathrm{O}^{+}+\mathrm{NO}_{3}^{-}$scores 1 overall <br> Correct species in square brackets can just score the arrow mark |


| Question |  |  | Answer |  | Mark | Guidance |
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| 1 | (e) | (ii) | $-\log (0.015) \checkmark=1.82(391) \checkmark$ |  | 2 | ALLOW 1.8 or more sf. ALLOW 'Ig' Answer alone scores 2 marks No ecf from first mark |
| 1 | (f) | (i) | $\begin{aligned} & \mathrm{H}_{2} \mathrm{SO}_{4}+\mathrm{HNO}_{3} \rightarrow \mathrm{NO}_{2}^{+}+\mathrm{H}_{2} \mathrm{O}+\mathrm{HSO}_{4}^{-} \\ & \mathrm{NO}_{2}^{+} \checkmark \text { rest correct } \checkmark \end{aligned}$ |  | 2 | $\text { ALLOW } 2 \mathrm{H}_{2} \mathrm{SO}_{4}+\mathrm{HNO}_{3} \rightarrow \mathrm{NO}_{2}^{+}+\mathrm{H}_{3} \mathrm{O}^{+}+2 \mathrm{HSO}_{4}^{-}$ <br> 'rest correct' means all other species correct with ' $\mathrm{NO}_{2}$ ' (incorrect sign or no sign) <br> IGNORE state symbols |
| 1 | (f) | (ii) | hydrogensulfate(VI) $\checkmark$ |  | 1 | Mark separately from f(i) IGNORE formula ALLOW 'hydrogen sulfate' and 'hydrogensulfate’ (i.e. no '(VI)') ALLOW ‘sulphate' replacing 'sulfate’ |
| 1 | (f) | (iii) | nitration OR nitrating (benzene/aromatics) OR making nitrobenzene $\checkmark$ |  | 1 | ALLOW 'electrophilic substitution of arene' |
|  |  |  |  | Total | 25 |  |


| Question |  |  | Answer | Mark | Guidance |
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| 2 | (a) | (i) | $\mathrm{CO}_{2} /$ carbon dioxide $\checkmark$ | 1 |  |
| 2 | (a) | (ii) | 2-hydroxypropanoic acid $\checkmark$ | 1 | ALLOW errors in gaps, commas, dashes ALLOW '2-hydroxylpropanoic acid |
| 2 | (b) | (i) | one correct 3d representation shown <br> two correct mirror images of 3d structures and either line between them or the words 'mirror images' | 2 | ALLOW groups on any carbons ambiguous attachments (eg COOH connected through O) <br> structures drawn as (but where two lines in the plane are shown (eg left-hand diagram here), they must not be at 90 or 180 degrees to each other) <br> ALLOW .."-י"'lll for $\qquad$ <br> ALLOW skeletal (or mixed) formulae <br> ALLOW structures breaking the 'two lines in the plane' rule above for second mark <br> NOT flat structures |
| 2 | (b) | (ii) | 1/ one $\checkmark$ | 1 |  |
| 2 | (c) | (i) | it has two: (carboxylic) acid / carboxyl groups / COOH groups $\checkmark$ | 1 | ALLOW 'dicarboxylic' and 'diprotic' |
| 2 | (c) | (ii) | lactic acid/it has a higher $\mathrm{pK}_{\mathrm{a}}$ value / lower $K_{\mathrm{a}} / 3.86$ is higher (ORA for malic acid) <br> higher pH / lower concentration (or fewer) $\underline{\underline{H}}^{ \pm}$ions / less dissociation into $\underline{H}^{ \pm}$ORA $\checkmark$ | 2 | Mark separately IGNORE references to 5.13/larger value for malic acid IGNORE 'less acidic' (which is in the stem) must refer to $\mathrm{H}^{+}$or pH |


| Question |  |  | Answer | Mark | Guidance |
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| 2 | (d) | (i) | $\mathrm{C}_{3} \mathrm{H}_{6} \mathrm{O}_{3} \rightleftharpoons \mathrm{C}_{3} \mathrm{H}_{5} \mathrm{O}_{3}^{-}+\mathrm{H}^{+} \checkmark$ | 1 | ALLOW $\mathrm{C}_{3} \mathrm{H}_{6} \mathrm{O}_{3}+\mathrm{H}_{2} \mathrm{O} \rightleftharpoons \mathrm{C}_{3} \mathrm{H}_{5} \mathrm{O}_{3}^{-}+\mathrm{H}_{3} \mathrm{O}^{+}$ $\mathrm{ORC}_{3} \mathrm{H}_{6} \mathrm{O}_{3}+\mathrm{aq}=\mathrm{C}_{3} \mathrm{H}_{5} \mathrm{O}_{3}^{-}+\mathrm{H}^{+}$ IGNORE lactic acid shown as a structural formula IGNORE state symbols watch out for incorrect anion formula |
| 2 | (d) | (ii) | $\begin{aligned} & K_{\mathrm{a}}=10^{-3.86} / \mathrm{pK}_{\mathrm{a}}=-\log \mathrm{K}_{\mathrm{a}} \\ & \mathrm{OR}-\log 1.4 \times 10^{-4}=3.86 \end{aligned}$ | 1 | No marks obviously for $1.4 \times 10^{-4}$ ALLOW 'inv(erse) $\log -3.86$ ' or ${ }^{\prime} \log ^{-1}(-3.86)$ ' ALLOW ' Ig ' for 'log' |
| 2 | (d) | (iii) | $\begin{aligned} & {\left[\mathrm{H}^{+}\right]=\sqrt{ }\left(1.4 \times 10^{-4} \times 0.1\right)\left(=3.74 \times 10^{-3}\right)^{\checkmark}} \\ & \mathrm{pH}=2.43 \checkmark \end{aligned}$ | 2 | First mark is for working* or correct evaluation Allow ecf for second mark* (if $\left[\mathrm{H}^{+}\right]$is smaller than $1 \times 10^{-2}$ ) ALLOW 2.4 or more sf Correct pH value alone scores two marks * provided working shows ' $\mathrm{H}^{+}=. .$. ' or ' $\left[\mathrm{H}^{+}\right]=. .$. ' |
| 2 | (d) | (iv) | $\begin{aligned} & {[\text { salt }] /[\text { acid }]=K_{\mathrm{a}} /\left[\mathrm{H}^{+}\right]} \\ & \text {OR [salt] }][\text { acid }]=1.4 \times 10^{-4} /(1 \mathrm{x}) 10^{-3} \checkmark \\ & =1.4 \times 10^{-4} / 1 \times 10^{-3}=0.14 \checkmark \end{aligned}$ | 2 | Correct answer (to 2 or more sf) scores both marks without working. Allow ecf for second mark for inverse only (7.1429 [to 2 or more sf] or $50 / 7$ ). This scores one mark even if no working. <br> ALLOW 1/7.1 OR 7/50 |
| 2 | (e) |  | C=O present in ester $\checkmark$ no $\mathrm{OH} /$ alcohol (groups) / no alkene $\checkmark$ | 3 | Both $\mathrm{C}=\mathrm{O}$ and ester must be mentioned <br> IGNORE references to other groups any unambiguous indication of the structure scores <br> Any additional incorrect structures are CON <br> NOT $\mathrm{CH}_{2}=\mathrm{CHCOOCH}\left(\mathrm{CH}_{3}\right) \mathrm{COOH}$ since alkene $\mathrm{C}-\mathrm{H}$ is 3000 up NOT ring with ether and anhydride as it is not an ester |


| Question |  |  | Answer | Mark | Guidance |
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| 2 | (f) |  |  <br> acid anhydride $\checkmark$ rest $\checkmark$ | 2 | Any unambiguous indication of the structure scores marks <br> Any additional incorrect structures are CON ALLOW any acid anhydride or just the anhydride group for the first mark <br> ALLOW (-)CO-O-CO(-) for anhydride part of formula |
| 2 | (g) | (i) | ethanal $\checkmark$ | 1 |  |
| 2 | (g) | (ii) |  | 1 | Any unambiguous representation of structure scores mark allow ambiguous attachment of any group <br> ALLOW anion formed at OH group $/ \mathrm{O}^{-} \mathrm{Na}^{+}$(or any Group I cation) |
| 2 | (g) | (iii) | nucleophilic $\checkmark$ addition $\checkmark$ | 2 | ALLOW any unambiguous indication of the words (eg circling) Mark separately <br> Any extra words CON correct answers |
| 2 | (g) | (iv) | ammonium ion / NH4 ${ }^{+} \checkmark \checkmark$ | 2 | ammonia/ $\mathrm{NH}_{3}$ scores one mark |
| 2 | (h) | (i) | orange/yellow to blue/green $\checkmark$ $\mathrm{CH}_{3} \mathrm{COOH} \checkmark$ | 2 | ALLOW any combinations of these colours but no others should be mentioned ALLOW a more displayed or skeletal formula |
| 2 | (h) | (ii) | $\mathrm{OH} /$ (secondary) alcohol / hydroxyl (group) (in lactic acid) <br> (lactic acid) reacts with reagent/acidified (potassium) dichromate gives same reaction / same colour change / same result / can be oxidised $\checkmark$ | 2 | 'primary/tertiary alcohol' is CON to first mark mark separately <br> ALLOW 'both have a colour change' |
|  |  |  | Total | 29 |  |


| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (a) |  | propane-1,2-diol $\checkmark$ | 1 | ALLOW errors in dashes, gaps and commas ACCEPT propan-1,2-diol |
| 3 | (b) | (i) | co-product is formed in the reaction that produces the (main) product <br> OR co-product is another substance (AW) that appears in the main reaction/equation $\checkmark$ <br> by-product is formed by side/other/unwanted reactions $\checkmark$ | 2 | must imply that co-product is another product of the reaction IGNORE 'process' (instead of 'reaction') IGNORE usefulness in either mark |
| 3 | (b) | (ii) | hydrolysis $\checkmark$ | 1 | ALLOW ‘saponification’ |
| 3 | (c) |  | water is only (other) product / water and no other product OR <br> glycerin is made from a renewable/carbon-neutral/ sustainable substance <br> OR <br> no toxic products $\checkmark$ | 1 | must imply that it is only water <br> IGNORE 'as a by-product' <br> ALLOW 'high atom economy' (but IGNORE 100\%) <br> IGNORE references to energy consumption IGNORE harmful/ 'no pollutants' |
| 3 | (d) |  | increases rate of reaction / faster reaction (AW) $\checkmark$ more frequent collisions (AW) $\checkmark$ | 2 | IGNORE references to equilibrium IGNORE what is colliding NOT just 'more collisions' IGNORE 'chance of collisions' |
| 3 | (e) | (i) |  | 1 | Arrow must start and end on bonds (or do so if the lines are continued with the same curvature) <br> Any other arrows on the glycerin structure are 'CON' |


| Question |  |  | Answer | Mark | Guidance |
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| 3 | (e) | (ii) | elimination $\checkmark$ | 1 |  |
| 3 | (e) | (iii) | (primary) alcohol / hydroxyl $\checkmark$ ketone / carbonyl $\checkmark$ | 2 | 'secondary' or 'tertiary' CON the alcohol mark Any extra groups CON a correct answer |
| 3 | (e) | (iv) | a IR: acetol has peak at $1705-1725 \checkmark$ <br> b $\mathrm{C}=\mathrm{O} /$ ketone $\checkmark$ <br> c NMR: acetol has peaks of height/area/ratio 3:2:1 (any order) <br> OR acetol has a peak at $2.0-2.7$ <br> OR two from the following: <br> - acetol has 3 peaks <br> - glycerin has 4 peaks <br> - glycerin has a doublet or triplet at 3.3-4.8 $\checkmark$ <br> d linking no. of peaks with proton/hydrogen environments in either compound $\checkmark$ <br> e all singlets/single peaks/unsplit/splitting of $1 \checkmark$ <br> f the neighbouring carbons have no hydrogens/protons | 6 | Please place ticks where points scored check it is the correct numerical range <br> IGNORE references to other peaks <br> c must mention peaks for this mark. <br> IGNORE any extra material even if wrong <br> d (QWC mark linking environments with no. of peaks) <br> ALLOW if hydrogen environments described rather than number given (eg acetol has $\mathrm{CH}_{3}, \mathrm{CH}_{2}$ and OH ) <br> Number of peaks must be stated for either compound and be the same as the number of environments for this mark to be awarded <br> e must imply all peaks are unsplit <br> f must actually say 'no hydrogens' not just state the rule references to any split peaks loses e but $\mathbf{f}$ can be scored if it is made for any peak |
| 3 | (f) | (i) | $+3 \checkmark$ | 1 | NOT 3+ |


| Question |  |  | Answer | Mark | Guidance |
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| 3 | (f) | (ii) | no effect $\checkmark$ <br> catalyst affects the rate of/speeds up forward and back reactions (equally) <br> OR affects/speeds up rate of attainment of equilibrium (AW) $\checkmark$ | 2 | ALLOW 'only temperature affects $K_{c}$ ' ALLOW 'provides a route of lower activation enthalpy/energy' IGNORE 'affects rate' without qualification mark separately |
| 3 | (g) |  | $\begin{aligned} & \text { calculation of both } M_{r} \text { values correctly } \checkmark \text { (glycerin } 92 \text {, } \\ & \text { prop glycol } 76) \\ & \%=(9 / 76) \times(92 / 15) \times 100=73 \% \checkmark \end{aligned}$ | 2 | ALLOW 92.0 and 76.0 <br> Answer on its own scores marks without reference to working ALLOW two or more sig figs allow any answer between 72 and 73 (intermediate rounding) <br> ALLOW ecf from incorrect $M_{\mathrm{r}}$ values (if clearly indicated as such or by working) |
|  |  |  | Total | 22 |  |


| Question |  |  | Answer |  |  |  | Mark | Guidance |
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| 4 | (a) | (i) |  |  |  |  | 2 | ALLOW - $\mathrm{CH}_{3}$ <br> ALLOW H atoms on benzene carbons IGNORE anything written over arrow |
| 4 | (a) | (ii) | ethanoyl chloride $\checkmark$ |  |  |  | 1 | ALLOW without gap between words |
| 4 | (a) | (iii) | electrophilic $\checkmark$ substitution $\checkmark$ |  |  |  | 2 | Mark separately <br> ALLOW mis-spellings that sound like 'electrophilic' IGNORE 'acylation / acetylation / ethanoylation Other words are CON to correct marks |
| 4 | (b) |  | Aluminium (compounds) are toxic $\mathbf{O R ~ H C l}$ is a toxic/corrosive/acidic (gas) / lowers pH of rivers/lakes (etc) $\checkmark$ |  |  |  | 1 | must name a product to score ALLOW aluminium chloride giving HCl with consequences of this |
| 4 | (c) |  | $\begin{array}{\|l} \hline\left(\mathrm{C}_{2} \mathrm{H}_{5}\right)_{4} \mathrm{~N}^{+} \\ \hline \mathrm{PF}_{6}^{-} \\ \hline \checkmark \checkmark \text { for eac } \\ \hline \end{array}$ | 4 <br> 6 | tetrahedral octahedral correct, one if | $\begin{aligned} & 107-110 \\ & 90 \\ & \hline \end{aligned}$ <br> one error | 4 | No ecf ALLOW 'tetrahedron' and/or 'octahedron' |
| 4 | (c) | (ii) | small/weak (AW): <br> electrostatic attraction / <br> attraction between ions / <br> force between ions / <br> ion-ion bonds/ionic bonds $\checkmark$ <br> small amount of energy required: <br> to separate (ions) / <br> break/overcome forces of attraction (or bonds) $\checkmark$ |  |  |  | 2 | IGNORE 'intermolecular', named imb, 'ion-dipole’ for first mark IGNORE answers that discuss removing electrons from ions first mark is for describing bonds between ions and implying they are weak <br> ALLOW breaking of any imb for second mark, provided energy is mentioned. second mark is for saying that little energy is required to break attraction/bonds AW <br> mark separately |


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| 4 | (d) | (i) | $\mathrm{CHCl}_{3}$ or more displayed $\checkmark$ | 1 | ALLOW $\mathrm{CHBr}_{3}$ <br> ALLOW triphenylmethane structure with one or two benzene rings replaced by Cl or $\mathrm{Br}\left(\mathrm{eg} \mathrm{CHCl}\left(\mathrm{C}_{6} \mathrm{H}_{5}\right)_{2}\right.$ or $\mathrm{C}_{13} \mathrm{H}_{11} \mathrm{Cl}$ or $\left.\mathrm{C}_{7} \mathrm{H}_{6} \mathrm{Cl}_{2}\right)$ |
| 4 | (d) | (ii) | a electrons excited/move up $\checkmark$ <br> b to (higher) energy level $\checkmark$ electron movement <br> c difference in levels related to frequency / $\Delta \mathrm{E}=\mathrm{hv} \checkmark$ relation between energy and frequency <br> d $\Delta$ E/energy gap or frequency is lower the greater the delocalisation / larger the chromophore ORA <br> OR comparison of size of energy gap (eg 'bigger gap for benzene' or 'big gap for benzene and small gap for dye') $\checkmark$ <br> comparison of benzene and dyes in terms of $\Delta E$ <br> e more delocalisation/larger chromophore for dyes than benzene ORA $\checkmark$ <br> reason: difference in delocalisation <br> f uv has: high frequency/low wavelength/more energy (than visible) ORA <br> OR benzene absorbs higher frequency than visible ORA $\checkmark$ <br> difference between uv and visible | 6 | Please place ticks where points scored <br> 'electrons move to higher energy levels' scores both $\mathbf{a}$ and $\mathbf{b}$ <br> QWC: only award point $\mathbf{b}$ if marking point $\mathbf{a}$ has been scored <br> IGNORE references to d-orbitals and ligands <br> c ALLOW $E=$ hv but only in context of energy gaps ALLOW ' $f$ ' for ' $v$ ' in $E=h v$ <br> d ALLOW in terms of excitation energies (eg 'it takes more energy to excite benzene's electrons') <br> ALLOW 'conjugated system' for 'chromophore/delocalisation' in d and $e$ <br> e ALLOW comment just for benzene or dyes, provided it is a comparison (eg 'benzene has less delocalisation'). Award for eg 'dyes big delocalisation, benzene small' in different parts of answer <br> If electrons dropping down and emission of light mentioned, max 2 out of 6 <br> Please place cross where 'emission' (AW) is mentioned |


| Question |  | Answer | Mark | Guidance |
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| 4 | (e) | $\mathbf{a}$ (Kekule structure would form) $\mathrm{C}_{6} \mathrm{H}_{6} \mathrm{Br}_{2}$ or $\mathrm{C}_{6} \mathrm{H}_{6} \mathrm{Br}_{4}$ or $\mathrm{C}_{6} \mathrm{H}_{6} \mathrm{Br}_{6}$ or their structures <br> b (Actual product is) $\mathrm{C}_{6} \mathrm{H}_{5} \mathrm{Br}$ or its structure <br> c substitution AND addition mentioned correctly <br> d benzene/delocalised structure is stable / addition removes stability $\checkmark$ <br> e benzene has delocalisation/ has delocalised electrons $\checkmark$ | 5 | Please place ticks where points scored <br> a and b ACCEPT structures without specific reference to where they come from <br> ALLOW skeletal structures <br> ALLOW correct names (1,2-dibromocyclohexa-(3,5)-diene, (1,2,3,4,5,6-)hexabromocyclohexane; bromobenzene) but IGNORE names if structures given <br> ALLOW full descriptions (eg 'one bromine added on each end of one double bond') <br> c IGNORE type of addition/substitution <br> e 'delocalisation'/'delocalised' must be spelled correctly at least once (QWC) |
|  |  | Total | 24 |  |


| Question |  |  | Answer | Mark | Guidance |
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| 5 | (a) |  | $3 \mathrm{~d}^{10} 4 \mathrm{~s}^{2} \checkmark$ | 1 | ALLOW either order ALLOW upper case letters but numbers must be superscripts |
| 5 | (b) | (i) | $\mathrm{Zn}^{2+}(\mathrm{g}) \rightarrow \mathrm{Zn}^{3+}(\mathrm{g})+\mathrm{e}^{(-)} \checkmark$ | 1 | IGNORE any state symbol on the electron ALLOW $\mathrm{Zn}^{2+}(\mathrm{g})-\mathrm{e}^{(-)} \rightarrow \mathrm{Zn}^{3+}(\mathrm{g})$ Zn symbol must be correct to score |
| 5 | (b) | (ii) | $\mathrm{d}^{10} /$ full (3)d $\checkmark$ | 1 | 'd' must be stated |
| 5 | (b) | (iii) | $\mathrm{ZnSO}_{4}{ }^{\checkmark}$ | 1 |  |
| 5 | (c) |  | $\mathrm{ZnCl}_{2} \cdot 2 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{ZnCl}(\mathrm{OH})+\mathrm{HCl}+\mathrm{H}_{2} \mathrm{O} \checkmark$ | 1 | IGNORE state symbols <br> ALLOW $\mathrm{Zn}(\mathrm{OH}) \mathrm{Cl}$ or either without brackets (but must be OH not HO) <br> ALLOW the zinc salt product hydrated with one water (correctly represented as ' $\mathrm{H}_{2} \mathrm{O}$ ') eg ' $\mathrm{ZnCl}(\mathrm{OH}) \cdot \mathrm{H}_{2} \mathrm{O}$ ' <br> ALLOW $\mathrm{ZnCl}^{+}+\mathrm{OH}^{-}$for salt |
| 5 | (d) | (i) | lattice enthalpy/energy (of zinc chloride) $\checkmark$ | 1 | ALLOW 'lattice formation energy/enthalpy' ALLOW 'enthalpy change of lattice (energy/enthalpy)', 'lattic' for 'lattice' <br> IGNORE $\Delta H_{\text {LE }}$ etc <br> IGNORE signs |
| 5 | (d) | (ii) | (Sum of enthalpy changes of hydration) $=-2772$ <br> Expression: <br> (Sum of enthalpy changes of hydration) $+2734 \checkmark$ <br> evaluation of answer ( -38 ) (with ecf) $\checkmark$ | 3 | correctly evaluated look to see if marked on diagram <br> expression or correct evaluation <br> ALLOW -(-2734) <br> -38 scores three overall*; <br> (+)38 scores two overall* <br> +325 scores two overall* <br> -5506 scores two overall** <br> +685 scores two overall* <br> 325 or -325 score one overall* <br> * whether any working shown or not |


| Question |  |  | Answer | Mark | Guidance |
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| 5 | (d) | (iii) | broken: hydrogen bonds (in water) ionic bonds / electrostatic forces / ion-ion bonds (in lattice) <br> made: ion-(permanent) dipole bonds/forces $\checkmark$ | 3 | Extra wrong bonds mentioned are CON to correct bonds IGNORE the species linked by the bonds (eg 'Zn and Cl') |
| 5 | (e) |  | a Entropy is a measure of disorder OR entropy (implies) more ways of arrangement OR entropy measures number of ways of arrangement <br> b (Particles in) solution (often) more disordered / have more ways of arrangement (than solids) <br> c Ca ions: <br> have large numbers of/more water molecules clustered round/ <br> OR attract large numbers of/more water molecules <br> OR are greatly/strongly hydrated <br> OR have strong attraction to water molecules $\checkmark$ <br> d Entropy loss of water / more order in water $\checkmark$ | 4 | Please place ticks where points scored <br> a If 'disorder' or 'ways of arrangement' are qualified it must be by plural 'particles', 'ions', 'molecules', or '(quanta of) energy' NOT singular 'particle' etc, 'compound', 'atoms' or 'element' which CON correct words <br> IGNORE 'substance’ <br> b IGNORE 'liquid' for 'solution'; must imply dissolving ALLOW 'molecules' or 'ions' or 'solute' for 'particles' NOT just 'higher entropy' <br> c 'Ca' can be implied eg 'small highly charged ions' must say 'many' water molecules or 'strongly' hydrated <br> d must refer to water |
| 5 | (f) |  | a (Solubility) increase <br> b $\Delta \mathrm{S}_{\text {tot }}=\Delta \mathrm{S}_{\text {sys }}-\Delta \mathrm{H} / \mathrm{T} \checkmark$ <br> c $\Delta \mathrm{H} / \mathrm{T}$ gets smaller as $T$ gets bigger (AW) $\checkmark$ <br> d $\Delta \mathrm{S}_{\text {tot }}$ gets more positive / less negative / increases $\checkmark$ | 4 | Please place ticks where points scored <br> b ALLOW $\Delta \mathrm{S}_{\text {tot }}=\Delta \mathrm{S}_{\text {sys }}+\Delta \mathrm{S}_{\text {surr }}$ and $\Delta \mathrm{S}_{\text {surr }}=-\Delta \mathrm{H} / T$ look out for this in any part of the answer <br> c ALLOW ' $-\Delta H / T / \Delta S_{\text {surr }}$ gets less negative'/ ' $\Delta \mathrm{S}_{\text {surr }}$ increases'/ <br> ' $-\Delta \mathrm{S}_{\text {surr }}$ decreases' <br> IGNORE ' $\Delta \mathrm{S}_{\text {surr }}$ decreases' ' $-\Delta \mathrm{H} / \mathrm{T}$ decreases' <br> d ALLOW 'total entropy gets...' IGNORE just 'entropy gets...' |
|  |  |  | Total | 20 |  |

