# Markscheme 

May 2018

## Chemistry

Higher level

## Paper 3

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## Section A

| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1. | a | i | consists of single/one sheet/layer «of carbon atoms» $\checkmark$ <br> graphene has no density measurement <br> OR <br> graphene has no distance between layers data <br> OR <br> graphene has large specific surface area «compared to graphite» $\checkmark$ | Do not accept "sp"" alone without reference to single/one sheet/layer. <br> Accept "thickness of one atom" OR "consists of a plane" for M1. | 2 |
| 1. | a | ii | Any one of these alternatives: <br> ALTERNATIVE 1 $\left\lvert\, \begin{aligned} & « \frac{1.3 \times 10^{11}}{76 \times 10^{6}} » \\ & 1.7 \times 10^{3} / 1711 \checkmark \end{aligned}\right.$ <br> ALTERNATIVE 2 <br> $1600 \times 76 \times 10^{6}=1.2 \times 10^{11}$ «is less than tensile strength of graphene» $\checkmark$ <br> ALTERNATIVE 3 <br> $\frac{1.3 \times 10^{11}}{1600}=8.1 \times 10^{7}$ «is greater than upper end of tensile strength for graphite» $\checkmark$ | Accept any value in the range 1700-27 083. Answer may be expressed in scientific notation or otherwise. <br> Accept any value calculated which is less than the graphene tensile strength based on a value chosen from within the $4.8-76 \times 10^{6}$ range. | 1 |

(Question 1a continued)

| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1. | a | iii | «graphene has a high electron mobility of» 15000-200000 «cm ${ }^{2} \mathrm{~V}^{-1} \mathrm{~s}^{-1}$ » $\checkmark$ | A specific value or range of values must be given. <br> Accept any value in the 15000-200000 «cm ${ }^{2} V^{-1} s^{-1} »$ range. | 1 |
| 1. | b |  | smaller/zero $\checkmark$ <br> no delocalized electrons/electrons are bound/electrons not free to move/electrons not free to roam <br> OR <br> localized electrons «in sigma bonds» <br> OR <br> large band gap $\checkmark$ | Accept "diamond is a dielectric" OR "diamond does not conduct electricity" for M2. <br> Award [1 max] for just "immobile/less mobile". <br> Award [2] for "electrons immobile «in diamond" due to the large band gap" OR "electrons «in diamond» immobile since electrons are localized «in the sigma bonds»". | 2 |


| Question |  | Answers | Total |  |  |
| :--- | :--- | :--- | :--- | :--- | :--- |
| 1. |  | shorter bonds in graphene <br> OR <br> bonds in graphene intermediate between single and double <br> OR <br> bond order in graphene is 1.33 <br> OR <br> delocalization creates stronger bonds <br> OR <br> shorter bonds are stronger $\checkmark$ <br> stronger/shorter bonds require higher temperature/faster thermal motion to be <br> altered <br> OR <br> stronger/shorter bonds require greater energy to be broken $\checkmark$ | $\mathbf{2}$ |  |  |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 2. | a |  | Any two of: <br> Ethene: «carbon-carbon» double bond AND Ethane: «carbon-carbon» single bond $\checkmark$ <br> ethene has a shorter carbon-carbon bond «than ethane» $\checkmark$ <br> Ethene: planar/two-dimensional/2-D AND Ethane: tetrahedral «carbons»/ three-dimensional/3-D <br> OR <br> Ethene: each carbon surrounded by three electron domains AND Ethane: each carbon surrounded by four electron domains <br> OR <br> different molecular geometries/shapes $\checkmark$ <br> rotation about carbon-carbon inhibited/blocked in ethene AND not in ethane $\checkmark$ <br> «H-C-C/H-C-H» bond angles different <br> OR <br> Ethene: «bond angles approximately» $120^{\circ}$ AND Ethane: $109.5 / 109^{\circ} \checkmark$ | Do not accept "different number of atoms/hydrogens/bonds" etc. <br> Accept "Ethene: unsaturated AND Ethane: saturated" OR "Ethene: has a double bond AND Ethane: does not" OR "Ethene: two flexible bonds between carbon atoms AND Ethane: one". <br> Accept any reasonable physical description of the two different molecular models based on a variety of kits for M1. <br> For ethene, accept any bond angle in the range 117-1220. <br> Award [2] if any two of the concepts listed are shown in a correctly labelled or annotated diagram. <br> Award [1 max] for two correct statements for either molecule but with no comparison given to the other. <br> Award [1 max] for suitable unlabeled diagrams of both compounds. | 2 max |


| Question |  |  | Answers | Notes | Total |
| :--- | :--- | :--- | :--- | :--- | :--- |
| 2. | b | $\mathbf{i}$ | 6 carbon atoms labelled in correct positions $\checkmark$ <br> both nitrogen atoms labelled in correct positions $\checkmark$ <br> bromine $\boldsymbol{A N D}$ chlorine atoms labelled in correct positions $\checkmark$ |  |  |

(Question 2b continued)

| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 2. | b | ii | accurate bond angles/lengths can be measured <br> OR <br> «using mathematical functions» can calculate expected shapes based on energy minimizations <br> OR <br> better visualization of possible bond rotations/conformation/modes of vibration <br> OR <br> can visualize macromolecules/proteins/DNA <br> OR <br> hydrogen bonding «networks» can be generated/allows intermolecular forces «of attraction» to be simulated <br> OR <br> more variety of visualization representations/can observe space filling <br> OR <br> can produce an electron density map/electrostatic potential map <br> OR <br> once model is generated file can be saved for future use/computer models can be shared globally by scientists <br> OR <br> helps design molecules of biological significance/assists in drug design «using libraries» <br> OR <br> can predict molecular interactions with solvents/can predict physical properties/can predict spectral data/can examine crystal structures <br> OR <br> «often» easier to construct/modify «model» $\downarrow$ | Accept "precise" for "accurate". <br> Accept "computer generated structural representation is normally what is expected in order to be published «in a scientific journal»". <br> Accept "easier to see different sizes of atoms/atomic radii". | 1 |

(Question 2b continued)

| Question |  | Answers | Notes |
| :--- | :--- | :--- | :--- | :--- | :--- |
| 2. | $\mathbf{b}$ | iii | bonds within ring have resonance <br> OR <br> contains delocalized «conjugated pi» electrons in ring $\boldsymbol{V}$ |
| There must be reference to a ring or <br> cyclic structure. |  |  |  |
| Accept "alternating single and double |  |  |  |
| bonds in a ring". |  |  |  |
| Accept "ring which shows |  |  |  |
| resonance/delocalization". |  |  |  |
| Accept "follows Hückel/4n +2 rule". |  |  |  |
| Do not accept "contains one or more |  |  |  |
| benzene rings". |  |  |  |

## Section B

## Option A - Materials

| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 3. | a | Alloy: mixture of metal with other metals/non-metals <br> OR mixture of elements that retains the properties of a metal $\checkmark$ Composite: reinforcing phase embedded in matrix phase $\checkmark$ | Award [1 max] for implying "composites only have heterogeneous/nonhomogeneous compositions". | 2 |
| 3. | b | effective for yttrium «but less/not for nickel» $\checkmark$ points on nickel graph do not lie on $« y=x »$ line OR cannot be used for low concentrations of nickel OR concentration of nickel is lower than recorded value $\checkmark$ | Accept "ICP-OES is more accurate for lower yttrium concentrations than higher concentrations" for M1. <br> Accept [Ni] and [Y] for concentrations of nickel and yttrium. <br> Accept "detection limit for yttrium is lower than for nickel" for M2. <br> Award [1 max] for "more accurate for yttrium at lower concentrations AND nickel at higher concentrations". | 2 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3. | C | i | Graph 1: determines wavelength of maximum absorption/maximum intensity «for vanadium» $\downarrow$ <br> Graph 2: determines absorption of known concentrations «at that wavelength» OR estimates [V]/concentration in a sample using «the signal» intensity $\checkmark$ | Do not accept just "determines maximum wavelength/ $\lambda_{\max }$ " for M1. <br> Do not accept "calibration curve" for M2. | 2 |
| 3. | c | ii | $\begin{aligned} & « 14950=392.19 x+147.62 » \\ & x=37.74 « \mu \mathrm{~g} \mathrm{~kg}^{-1} » \end{aligned}$ | Answer must be given to four significant figures. <br> Do not accept values obtained directly from the graph. | 1 |
| 3. | c | iii | vanadium reduced in first reaction AND oxidized in second reaction <br> OR <br> $\mathrm{V}_{2} \mathrm{O}_{5}$ oxidizes $\mathrm{SO}_{2}$ in first reaction $\boldsymbol{A N D} \mathrm{VO}_{2}$ reduces $\mathrm{O}_{2}$ in second reaction <br> OR <br> vanadium returns to original oxidation state «after reaction» $\checkmark$ <br> provides an alternative reaction pathway/mechanism «with a lower activation energy» $\downarrow$ | Do not accept "reactants adsorb onto surface AND products desorb". <br> Accept "oxidation number" for "oxidation state". | 2 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4. | a | i | $2 \checkmark$ |  | 1 |
| 4. | a | ii | $n \lambda=2 d \sin \theta$ <br> OR $\begin{aligned} & \theta=\sin ^{-1}\left(\frac{n \lambda}{2 d}\right) \checkmark \\ & \theta=« \sin ^{-1}\left(\frac{150}{2 \times 303}\right)=» 14.3 «^{\circ} » \end{aligned}$ | Award [2] for correct final answer. | 2 |
| 4. | a | iii | $m=« \frac{50.94}{6.02 \times 10^{23}}=» 8.46 \times 10^{-23} « \mathrm{~g} » \downarrow$ |  | 1 |
| 4. | a | iv | $\begin{aligned} & « 303 \mathrm{pm}=303 \times 10^{-10} \mathrm{~cm} \\ & V=«\left(303 \times 10^{-10}\right)^{3}=» 2.78 \times 10^{-23} « \mathrm{~cm}^{3} » \end{aligned}$ |  | 1 |
| 4. | a | v | $\begin{aligned} & « 8.46 \times 10^{-23} \mathrm{~g} \times 2=» 1.69 \times 10^{-22} \text { «g» } \\ & d=« \frac{1.69 \times 10^{-22} \mathrm{~g}}{2.78 \times 10^{-23} \mathrm{~cm}^{3}}=» 6.08 « \mathrm{~g} \mathrm{~cm}^{-3} » \end{aligned}$ | Accept any value in the range $6.07-6.09 « \mathrm{~g} \mathrm{~cm}{ }^{-3} » .$ <br> Award [2] for correct final answer. | 2 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4. | b | i | Any one of these alternatives: <br> ALTERNATIVE 1 <br> disrupt enzyme binding sites $\checkmark$ <br> which can inhibit/over-stimulate enzymes $\checkmark$ <br> ALTERNATIVE 2 <br> disrupt endocrine system $\checkmark$ <br> because they compete for active sites of enzymes/cellular receptors $\checkmark$ <br> ALTERNATIVE 3 <br> form complexes/coordination compounds $\checkmark$ <br> which can bind to enzymes $\checkmark$ <br> ALTERNATIVE 4 <br> act as oxidizing/reducing agents <br> OR <br> act as catalysts $\checkmark$ <br> which can initiate unwanted reactions $\checkmark$ | Accept "can undergo oxidationreduction reactions" for M1 in Alternative 4. | 2 |
| 4. | b | ii | $\mathrm{V}^{4+}(\mathrm{aq})+\mathrm{H}_{2} \mathrm{O}_{2}(\mathrm{aq}) \rightarrow \mathrm{V}^{5+}(\mathrm{aq})+\mathrm{OH}^{-}(\mathrm{aq})+\bullet \mathrm{OH}(\mathrm{aq}) \checkmark$ | Do not accept • on H. <br> Accept answer without $\bullet$. | 1 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5. | a |  |  | Do not accept syndiotactic (alternating orientation of the $\mathrm{CH}_{3}$ groups), eg, <br> for M1 or M2. <br> Accept any correct atactic ordering of $\mathrm{CH}_{3}$ groups. <br> Penalize missing hydrogens or incorrect bond connectivities once only. <br> Accept skeletal structures. <br> Ignore continuation bonds, brackets and " $n$ " indices in structures. | 2 |
| 5. | b |  | Any two of: <br> Recycling: shredded/melted/reformed AND Reuse: used in its current form $\checkmark$ recycling is more energy intensive «than reusing» $\checkmark$ <br> recycling degrades the quality of plastic but reusing «typically» does not $\sqrt{ }$ <br> recycling breaks down original product to form a new product whereas reuse extends product life $\checkmark$ |  | 2 max |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5. | c | i | Any one of these alternatives: <br> ALTERNATIVE 1 <br> Polyester: produced by condensation/esterification polymerization $\checkmark$ <br> Polyethene: produced by addition polymerization $\checkmark$ <br> ALTERNATIVE 2 <br> Polyester: reaction between monomers/molecules containing two functional groups per molecule $\checkmark$ <br> Polyethene: reaction between monomers/molecules containing a carbon-carbon double bond/C=C $\checkmark$ <br> ALTERNATIVE 3 <br> polyester polymerization forms a by-product/ $/ \mathrm{H}_{2} \mathrm{O} \checkmark$ <br> polyethene has no by-products/100\% atom economy $\checkmark$ | Accept the names of different catalysts used for each polymerization as an alternative answer. | 2 |
| 5. | c | ii | more pliable/flexible materials <br> OR <br> more durable/non-corrosive/longer-lasting materials <br> OR <br> greater variety of materials <br> OR <br> lower density <br> OR <br> can be clear/translucent $\checkmark$ | Accept "more adaptable". <br> Do not accept just "more useful". | 1 |



## Option B — Biochemistry

| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 7. | a |  | Type of reaction: <br> condensation <br> OR <br> esterification/triesterification <br> OR <br> nucleophilic substitution/nucleophilic displacement/SN2 <br> By-product: <br> water/ $\mathrm{H}_{2} \mathrm{O} \checkmark$ | Do not accept just "substitution/displacement". | 2 |
| 7. | b |  | ALTERNATIVE 1 $\begin{aligned} & « \frac{334}{253.8}=» 1.32 \text { AND « } \frac{100}{304.5}=» 0.328 \\ & « \frac{1.32}{0.328} \approx » 4 \checkmark \end{aligned}$ <br> ALTERNATIVE 2 $\begin{aligned} & « 334 \times \frac{304.5}{100} \approx » 1017 \checkmark \\ & « \frac{1017}{253.8} \approx » 4 \checkmark \end{aligned}$ | Award [2] for correct final answer. | 2 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 7. | c |  |  <br> glycerol backbone as circled $\checkmark$ ester section as boxed $\checkmark$ | Accept a skeletal structure. <br> Penalize missing hydrogens or incorrect bond connectivities once only in Option B. <br> Accept condensed formula for ester. <br> Do not accept structures with one or two ester groups. | 2 |
| 7. | d |  | has affected consumption of trans-fats/cis-fats/saturated fats/unsaturated fats/ hydrogenated/artificially altered fats <br> OR <br> reduce/eliminate trans-fats/increase in cis-fats <br> OR <br> reduce/eliminate saturated fats <br> OR <br> increase unsaturated fats $\checkmark$ | Do not accept "decrease in fat" alone. <br> Accept "lipid" for "fats". | 1 |
| 7. | e |  |  | Ignore significant figures in M1. <br> Award [2] for correct final answer. <br> Award [1 max] for incorrect significant figures in final answer. | 2 |



| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 8. | d | ALTERNATIVE 1 $« \mathrm{pH}=6.36+\log \left(\frac{2.50 \times 10^{-2}}{1.25 \times 10^{-3}}\right)=»$ <br> $7.66 \checkmark$ <br> ALTERNATIVE 2 <br> $« K_{a}=4.4 \times 10^{-7}=\left[\mathrm{H}^{+}\right]\left(\frac{2.50 \times 10^{-2}}{1.25 \times 10^{-3}}\right),\left[\mathrm{H}^{+}\right]=2.2 \times 10^{-8} \mathrm{~mol} \mathrm{dm}^{-3}$ » «pH =» 7.66 V | Do not accept " $<\mathrm{pH}=$ » 8 ". | 1 |
| 8. | e |  | Penalize missing hydrogens or incorrect bond connectivities once only in Option B. <br> Wedges AND dashes must be used. | 1 |
| 8. | f | $\text { « } \frac{0.725}{49650 \mathrm{dm}^{3} \mathrm{~cm}^{-1} \mathrm{~mol}^{-1} \times 1.00 \mathrm{~cm}}=» 1.46 \times 10^{-5} \text { «mol dm}{ }^{-3} » \checkmark$ |  | 1 |
| 8. | g | 0.65 « $\mu \mathrm{g} \mathrm{cm}^{-3} \geqslant \checkmark$ | Accept any value in the range $0.60-0.70 « \mu \mathrm{~g} \mathrm{~cm}{ }^{-3}$ ». | 1 |


| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 9. |  | Any two of: replaces plastics with biodegradable/starch/cellulose based plastics $\checkmark$ <br> use enzymes instead of polluting detergents/phosphates <br> OR <br> use of enzymes means lower temperatures can be used <br> OR <br> use enzymes instead of emulsifiers to treat oil spills <br> OR <br> use enzymes to produce esters at lower temperatures/without sulfuric acid $\checkmark$ <br> replace organic/toxic solvents with carbon dioxide $\checkmark$ <br> replace polymers from fossil fuel with bamboo/renewable resources $\checkmark$ <br> develop paint resins reducing production of volatile compounds «when paint is applied» $\checkmark$ <br> industrial synthesis of ethanoic/acetic acid from methanol and carbon monoxide has $100 \%$ atom economy $\checkmark$ <br> energy recovery $\checkmark$ | Accept formulas for names. <br> Award mark for any other reasonable specific green chemistry example that prevents the release of pollutants/toxic chemicals into the environment by changing the method or the materials used. <br> Do not award mark for methods that involve clean-up of pollutants from the environment such as host-guest chemistry or alternative energy sources. | 2 max |


| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 10. | a | Vitamin A: <br> fat soluble/soluble in non-polar solvents AND non-polar/long hydrocarbon backbone/chain $\checkmark$ <br> Vitamin C: water soluble AND contains 4 hydroxyl groups/contains many hydroxyl groups/forms «many» H -bonds with water $\checkmark$ | Accept "Vitamin A: fat soluble/soluble in non-polar solvents as it contains only one hydroxyl group whose H-bonds with water are not strong enough to overcome London/dispersion/vdW forces between Vitamin A molecules". <br> Accept "lipid" for "fats". <br> Accept "alcohol" OR "hydroxy" OR "OH groups" for "hydroxyl" but not "hydroxide". <br> Award [1 max] for "Vitamin A: fat soluble AND Vitamin C: water soluble" with no or incomplete explanation. | 2 |
| 10. | b | vitamin A oxidized to «11-cis-»retinal $\downarrow$ <br> extended conjugation <br> OR <br> extensive delocalization $\checkmark$ <br> cis-retinal converts to trans-retinal through absorption of light $\checkmark$ | Accept "vitamin A/hydroxyl/hydroxy/alcohol/ $\mathrm{CH}_{2} \mathrm{OH}$ group oxidized to aldehyde/CHO «group in retinal»". | 3 |


| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 11. | a |  <br> curve below original curve «showing lower affinity for oxygen» beginning at $0 \checkmark$ | Award mark if end of student curve does not finish at same location as original curve. | 1 |
| 11. | b | Any two of: foetal hemoglobin has higher affinity for oxygen «than normal hemoglobin» $\checkmark$ foetal hemoglobin is less sensitive to inhibitors/2,3-bisphosphoglycerate/ 2,3-BPG/DPG «than normal hemoglobin» $\checkmark$ <br> foetal hemoglobin contains two gamma units instead of the two beta units found in adult hemoglobin $\checkmark$ |  | 2 max |


| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 12. |  | Any two of: pentose «sugar» <br> OR <br> deoxyribose $\checkmark$ <br> phosphate «group» $\checkmark$ <br> «organic» nitrogenous base <br> OR <br> nucleobase <br> OR <br> nucleic base <br> OR <br> purine <br> OR <br> pyrimidine $\checkmark$ | Accept names or formulas. <br> Accept "ribose" for M1. <br> Do not accept "phosphoric acid". <br> Accept the four bases together: "adenine, cytosine, thymine, guanine". | 2 max |

Option C - Energy

| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 13. | a |  | Any two of: high energy content/high energy density/high specific energy <br> OR high enthalpy of combustion/very exothermic enthalpy of combustion $\checkmark$ shortage of alternatives <br> OR <br> alternatives are expensive <br> OR <br> oil is relatively cheap <br> OR <br> oil is «still» abundant/common $\checkmark$ <br> well-established technology <br> OR <br> easy for consumers to obtain <br> OR <br> commonly used $\checkmark$ <br> easy to store <br> OR <br> easy to transport <br> OR <br> easy to extract $\checkmark$ <br> produces energy at a reasonable rate $\checkmark$ | Accept "high potential energy" for M1. | 2 max |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 13. | b | i | fuels can be compressed more without undergoing «unwanted» auto-ignition $\checkmark$ | Accept "burns smoother without undergoing «unwanted» auto-ignition" OR "fuel does not auto-ignite". | 1 |
| 13. | b | ii | produces more branched chain hydrocarbons «with higher octane rating» OR produces aromatics «which have higher octane rating» OR produces cyclohexanes «which have higher octane rating» $\checkmark$ | Accept "increase branches". <br> Do not accept "produces benzene". <br> Do not penalize for "benzene" if penalty applied in 2.b.iii. <br> Accept "produces cyclic structures". | 1 |
| 13. | C |  | $\begin{aligned} & n=6 \checkmark \\ & « \Delta G^{\ominus}=-n F E^{\ominus}=6 \mathrm{~mol} \times 96500 \mathrm{C} \mathrm{~mol}^{-1} \times 0.576 \mathrm{~V}=»-333504 \mathrm{~J} /-334 \mathrm{~kJ} \\ & \text { «Efficiency }=\frac{\Delta G}{\Delta H}=\frac{-334}{-726}=» 0.459 / 45.9 \% \text {, } \end{aligned}$ | Award [3] for correct final answer. | 3 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 14. | a |  | Any three of: <br> IR/long wavelength/low frequency radiation radiated/emitted by the Earth's «surface absorbed in the bonds» $\checkmark$ <br> bond length $/ \mathrm{C}=\mathrm{O}$ changes <br> OR <br> «asymmetric» stretching of bonds <br> OR <br> bond angle/OCO changes $\checkmark$ <br> polarity/dipole «moment» changes <br> OR <br> dipole «moment» created «when molecule absorbs IR» $\checkmark$ <br> «some of» energy is then re-radiated towards «the surface of the» Earth $\checkmark$ | Do not accept terms such as "reflect" OR "bounced" OR "trapped". | 3 max |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 14. | b |  | Any two of: <br> $\mathrm{H}_{2} \mathrm{O}$ AND «relatively» greater abundance/stable concentration/less effective at absorbing radiation/lower GWP so not much overall effect on global warming/climate change $\checkmark$ <br> $\mathrm{CH}_{4} / \mathrm{N}_{2} \mathrm{O} / \mathrm{CFCs} / \mathrm{SF}_{6} / \mathrm{O}_{3} / \mathrm{HCFCs}$ AND more effective «than $\mathrm{CO}_{2}$ » at absorbing radiation/higher GWP so could contribute to global warming/climate change $\checkmark$ <br> PFCs/SF $/{ }_{6} / \mathrm{NF}_{3} /$ Some CFCs $\boldsymbol{A N D}$ have very long life in atmosphere so could contribute «in the future» to global warming/climate change $\checkmark$ | Accept names or formulas. <br> Accept two different gases with the same effect for [2]. <br> Award [1 max] for identifying the names/formulas of two greenhouse gases. <br> Accept "greenhouse factor" for "GWP" but not just "greenhouse effect". <br> For M3, do not allow "CFC" alone as only some have long lifetimes (eg, CFC-115, CFC-113). | 2 max |


| 15. | a | « $\frac{813 \mathrm{~K}-296 \mathrm{~K}}{813 \mathrm{~K}} \times 100 »=64 « \% » \checkmark$ |  |  |
| :--- | :--- | :--- | :--- | :--- |
| 15. | $\mathbf{b}$ | $35 \%$ of chemical/potential energy available in coal is transformed to <br> electricity/electrical energy $\checkmark$ <br> not all chemical energy from burning fuel transferred into heating water <br> OR <br> energy dispersed elsewhere/energy lost due to friction of moving parts <br> OR <br> heat loss to the surroundings $\checkmark$ | Accept "stored energy" for <br> "potential energy". | 2 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 16. | a |  | Award [1] for one similarity: <br> both increase binding energy/energy yield «per nucleon» <br> OR <br> mass loss/defect in both «nuclear» reactions/mass converted to energy «from $E=m c^{2} »$ <br> OR <br> both produce ionizing radiation $\checkmark$ <br> Award [2 max] for any two differences: <br> in fusion, light nuclei combine to form heavier ones AND in fission, heavier nuclei split into lighter ones $\checkmark$ <br> fission produces radioactive/nuclear waste AND fusion does not $\checkmark$ <br> fission is caused by bombarding with a neutron «or by spontaneous fission» AND fusion does not <br> OR <br> fission can initiate a chain reaction $A N D$ fusion does not $\checkmark$ <br> fusion releases more energy per unit mass of fuel than fission $\checkmark$ <br> fuel is easier to obtain/cheaper for fusion reactions $\checkmark$ <br> fission reactions can be controlled in a power plant $\boldsymbol{A N D}$ fusion cannot «yet» $\checkmark$ <br> fusion reactor less likely to cause a large-scale technological disaster compared to fission $\checkmark$ <br> fusion less dangerous than fission as radioactive isotopes produced have short half-lives so only cause a threat for a relatively short period of time $\checkmark$ <br> fusion is in experimental development $\boldsymbol{A N D}$ fission used commercially $\checkmark$ | Accept "small nuclei" OR "smaller atomic masses of nuclei" for "light nuclei" AND "large nuclei" OR "greater atomic masses of nuclei" for "heavier nuclei". <br> Do not accept "no/less waste produced for fusion". <br> Accept "higher specific energy for fusion". | 3 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 16. | b |  | $\frac{1}{64} / \frac{1}{2^{6}} / 0.016 \checkmark$ | Accept "1.6\%". | 1 |
| 16. | C | i | $M_{r}\left({ }^{235} \mathrm{UF}_{6}\right)=235+(19.00 \times 6) / 349$ <br> OR | Award [2] for correct final answer. Do not accept "1.00" OR "0.996". | 2 |

(continued...)
(Question 16c continued)

| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 16. | c | ii | UF6: Structure: octahedral «solid»/square bipyramidal «solid»/«simple» molecular solid/simple molecule AND Bonding: covalent $\checkmark$ | Accept " $U_{6}$ : Structure: octahedral "solid»/square bipyramidal «solid»/«simple» molecular solid/simple molecule AND weak intermolecular/London/dispersion/van der Waals'/vdW forces". <br> Accept "non-polar molecule" for "«simple» molecular solid". |  |
|  |  |  | $\mathrm{UO}_{2}$ : Structure: crystal/lattice/network «solid»/«resembles» fluorite AND Bonding: «partly» covalent $\sqrt{ }$ | Accept "giant molecular" OR "macromolecular" for "network". <br> Accept "ionic/electrostatic attractions «between ions»" for bonding in $\mathrm{UO}_{2}$. <br> Award M2 for " $\mathrm{UO}_{2}$ : network covalent/covalent network/giant covalent" OR " $\mathrm{UO}_{2}$ : network ionic/giant ionic". | 3 |
|  |  |  | UF ${ }_{6}$ sublimes/evaporates/boils at low temperature $\checkmark$ | For M1 and M2 award [1 max] for two correct structures OR two bonding types. <br> Accept any specified low temperature in the range $56-65^{\circ} \mathrm{C}$. |  |


| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 17. | a | $\mathrm{C}_{7} \mathrm{H}_{15} \mathrm{COOC}_{5} \mathrm{H}_{11}(\mathrm{I})+\mathrm{CH}_{3} \mathrm{OH}(\mathrm{I}) \rightarrow \mathrm{C}_{7} \mathrm{H}_{15} \mathrm{COOCH}_{3}(\mathrm{I})+\mathrm{C}_{5} \mathrm{H}_{11} \mathrm{OH}(\mathrm{I})$ <br> OR $\mathrm{C}_{13} \mathrm{H}_{26} \mathrm{O}_{2}(\mathrm{I})+\mathrm{CH}_{4} \mathrm{O}(\mathrm{I}) \rightarrow \mathrm{C}_{9} \mathrm{H}_{18} \mathrm{O}_{2}(\mathrm{I})+\mathrm{C}_{5} \mathrm{H}_{12} \mathrm{O}(\mathrm{I})$ <br> OR | Accept correct equation in any format eg, skeletal, condensed structural formula, etc. <br> Accept equations with equilibrium arrow. | 1 |
| 17. | b | less viscous «and so does not need to be heated to flow» OR <br> less likely to undergo incomplete combustion <br> OR <br> fewer intermolecular/London/dispersion forces <br> OR <br> vaporizes easier $\checkmark$ | Ignore equation and products in 17 a. <br> Accept "van der Waals'/vdW" for "London". | 1 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 18. | a |  | ALTERNATIVE 1 <br> B/Ga in circle AND Type of semiconductor: p-type $\checkmark$ showing 3 electron pairs AND one lone electron «and hole» $\checkmark$ <br> ALTERNATIVE 2 <br> P/As in circle AND Type of semiconductor: n-type $\checkmark$ showing 4 electron pairs $\boldsymbol{A N D}$ one non-bonded electron $\checkmark$ | Accept any group 13 element labelled as p-type. <br> Accept showing 7 electrons. <br> Accept any group 15 element labelled as n-type. <br> Accept showing 9 electrons. <br> Accept dots or crosses for electrons. | 2 |
| 18. | b | i | conjugated $\mathrm{C}=\mathrm{C} /$ carbon-carbon double bonds <br> OR <br> «multiple» alternating $\mathrm{C}=\mathrm{C} /$ carbon-carbon double bonds <br> OR <br> «extensive electron» conjugation/delocalization <br> OR <br> «many» fused/conjugated aromatic/benzene rings $\checkmark$ |  | 1 |
| 18. | b | ii | complex B has greater conjugation/delocalization $\checkmark$ |  | 1 |

Option D - Medicinal chemistry

| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 19. |  | $L D_{50}$ : amount/dose that kills $50 \%$ of the population $\checkmark$ <br> $T D_{50}$ : amount/dose that negatively affects/produces toxic effects in $50 \%$ of the population $\checkmark$ | Award [1 max] for " $L D_{50}$ used in animal trials AND TD 50 used in human studies". | 2 |


| Question |  |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 20. | a | i | «irreversibly» binds/bonds to enzyme/transpeptidase <br> OR <br> inhibits enzyme/transpeptidase «in bacteria» that produces cell walls <br> OR <br> prevents cross-linking of bacterial cell walls $\checkmark$ <br> cells absorb water AND burst <br> OR <br> cells cannot reproduce $\checkmark$ |  | 2 |
| 20. | a | ii | modify side chain $\checkmark$ |  | 1 |
| 20. | b |  | condensation <br> OR <br> esterification <br> OR <br> nucleophilic substitution/nucleophilic displacement/S $\mathfrak{N} 2 \checkmark$ | Do not accept just "substitution/displacement". | 1 |


| Question |  | Answers | Notes |
| :--- | :--- | :--- | :--- | :--- | :--- |
| 20. | $\mathbf{c}$ | water causes hydrolysis <br> OR <br> aspirin reacts with water $\checkmark$ <br> heat increases the rate of hydrolysis <br> OR <br> heat increases the rate of the reaction with water $\checkmark$ | Accept "aspirin will convert into <br> salicylic/ethanoic acid". <br> Do not accept "aspirin dissolves in <br> water" OR "aspirin absorbs water/is <br> hygroscopic". |

21. $\quad$ morphine has hydroxyl/OH groups/is more polar AND diamorphine has ester/ethanoate/acetate groups/is less polar/is lipid soluble $\checkmark$
crossing blood brain barrier is easier for non-polar/less polar compounds/for lipid soluble compounds $\checkmark$

Accept "alcohol/hydroxy" for "hydroxyl" but not "hydroxide".

Accept "fats" for "lipid".

| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 22. | a | $2 \mathrm{HCl}(\mathrm{aq})+\mathrm{CaCO}_{3}(\mathrm{~s}) \rightarrow \mathrm{H}_{2} \mathrm{O}(\mathrm{I})+\mathrm{CO}_{2}(\mathrm{~g})+\mathrm{CaCl}_{2}(\mathrm{aq}) \checkmark$ | Accept ionic equation: $2 \mathrm{H}^{+}(\mathrm{aq})+\mathrm{CO}_{3}{ }^{2-}(\mathrm{aq}) \rightarrow \mathrm{CO}_{2}(\mathrm{~g})+\mathrm{H}_{2} \mathrm{O}(\mathrm{I})$ | 1 |
| 22. | b | « $\frac{0.750 \times 2}{100.09}=» 0.0150$ «mol HCl» |  | 1 |
| 22. | C | inhibits the secretion of stomach acid/ $/ \mathrm{H}^{+} \checkmark$ <br> «active metabolites» bind «irreversibly» to «receptors of the» proton pump $\downarrow$ | Do not accept "hydrogen/ $/ H^{\prime} / H_{2}$ " for " $\mathrm{H}^{+}$". <br> Accept "PPI/proton pump inhibitor" for M2. <br> Accept " $\mathrm{H}^{+} / K^{+}$ATPase" for "proton pump". | 2 |


| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 23. |  | Any two of: hydroxyl $\sqrt{ }$ carboxyl/carbonyl $\checkmark$ ether $\checkmark$ amido/carbonyl $\checkmark$ | Accept "alcohol/hydroxy" for "hydroxyl", "carboxylic acid" for "carboxyl" and "amide/carboxamide" for "amido". <br> Accept "amino/amine" OR "imine/imino" but these are not correct as they are part of the guanidino group. <br> Accept "alkenyl/alkene/carbon to carbon double bond" but not "C=C" OR "carbon double bond". <br> Accept "carbonyl" only once. <br> Accept "heterocyclic ring" for "ether". | 2 max |






| Question |  | Answers | Notes | Total |
| :---: | :---: | :---: | :---: | :---: |
| 27. | a | $\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OH}(\mathrm{g})+\mathrm{O}_{2}(\mathrm{~g}) \rightarrow \mathrm{CH}_{3} \mathrm{COOH}(\mathrm{aq})+\mathrm{H}_{2} \mathrm{O}(\mathrm{I}) \checkmark$ | Accept any correct formula for reactants and products. | 1 |
| 27. | b | R-OH: <br> 1.0-6.0 «ppm» AND 1 H $\mathrm{R}-\mathrm{O}-\mathrm{CH}_{2}-:$ <br> 3.3-3.7 «ppm» AND 2 H , $-\mathrm{CH}_{3}:$ <br> 0.9-1.0 «ppm» AND 3 H , | Award [1] for the ratio of 1:2:3 (in any order). <br> Award [2] for three correct chemical shifts without integration. <br> Award [1] for two correct chemical shifts without integration. <br> For each chemical shift accept a specific value within the range. <br> Assignment of proton to fragment (eg, $\mathrm{R}-\mathrm{OH}$ ) is not required in each case. | 3 |

