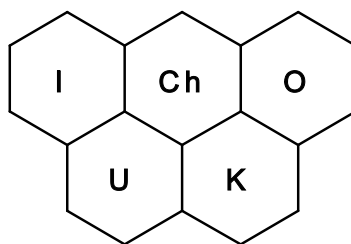




**INEOS**  
THE WORD FOR CHEMICALS



**48<sup>th</sup> INTERNATIONAL  
CHEMISTRY OLYMPIAD  
2016  
UK Round One  
MARK SCHEME**

Although we would encourage students to always quote answers to an appropriate number of significant figures, do not penalise students for significant figure errors. Allow where a student's answers differ slightly from the mark scheme due to the use of rounded/non-rounded data from an earlier part of the question.

In general error carried forward can be applied. We have tried to indicate where this may happen in the mark scheme.

For answers with missing or incorrect units, penalise one mark for the first occurrence in **each** question and write **UNIT** next to it. Do not penalise for subsequent occurrences in the same question.

Organic structures are shown in their skeletal form, but also accept displayed formulae as long as the representation is unambiguous. Benzene rings may be drawn with localised or delocalised bonding.

*Comments in blue have been added to some questions, these are not required for the marks, but may be of interest in subsequent discussion on the questions.*

Question	1	2	3	4	5	Total
Marks Available	12	23	31	21	13	100

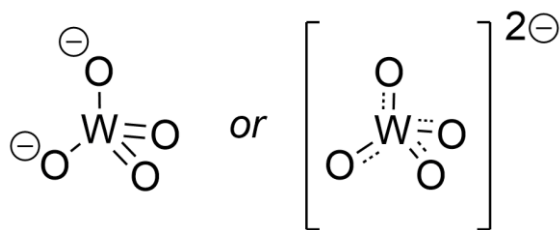
**1. This question is about energy storage using a chemical cycle**

- (a) (i)  $\Delta_r H^\circ (\text{reaction (3)}) = \Delta_f H^\circ (\text{I}_{2(\text{g})}) - 2 \Delta_f H^\circ (\text{HI}) = (62.4 - 2 \times 26.5) \text{ kJ mol}^{-1}$   
 $= +9.4 \text{ kJ mol}^{-1}$  1  
*Plus sign not required.*
- (ii)  $\Delta S^\circ (\text{reaction (3)}) = (131 + 261 - 2 \times 207) \text{ J K}^{-1} \text{ mol}^{-1}$   
 $= -22 \text{ J K}^{-1} \text{ mol}^{-1}$  1  
*Minus sign must be present.*
- (iii)  $\Delta_r G^\circ (\text{reaction (3)}) = 9.4 \text{ kJ mol}^{-1} - (298 \text{ K} \times -0.022 \text{ kJ K}^{-1} \text{ mol}^{-1})$   
 $= +16.0 \text{ kJ mol}^{-1}$  1  
*Plus sign not required. Allow error carried forward from (i) and/or (ii)*
- (iv)  $K_{298} = \exp(-\Delta G/RT) = \exp(15956 / (8.314 \times 298))$   
 $= 1.60 \times 10^3$  2  
*Allow error carried forward from (iii). Do not penalise if equilibrium constant has units.*
- (v)  $\Delta_r G^\circ (\text{reaction (3)}) = 9.4 \text{ kJ mol}^{-1} - (723 \text{ K} \times -0.022 \text{ kJ K}^{-1} \text{ mol}^{-1})$   
 $= 25.3 \text{ kJ mol}^{-1}$   
 $K_{723} = \exp(-\Delta G/RT) = \exp(25306 / (8.314 \times 723))$   
 $= 0.0148$  2  
*Allow error carried forward from (i) and/or (ii). Do not penalise if equilibrium constant has units.*
- (b) Products of reaction (1) cancel out when they occur in the following proportion:  
 $2 \times \text{reaction (1)} + \text{reaction (2)} + 2 \times \text{reaction (3)}.$   
 This simplifies down to the following reaction:  
 $2\text{H}_2\text{O} (\text{g}) \rightarrow 2\text{H}_2 (\text{g}) + \text{O}_2 (\text{g})$  2  
*State symbols not required. Accept the equation with mole ratio 1:1:½. Award 1 mark if the reactions are combined in the correct ratio but simplifying is done incorrectly.*
- (c)  $2 \times \Delta_r H^\circ (1) + \Delta_r H^\circ (2) + 2 \times \Delta_r H^\circ (3) = -2 \times \Delta_f H^\circ (\text{H}_2\text{O}_{(\text{g})})$   
 $[2 \times \Delta_r H^\circ (1) + 439 + 2 \times 9.4] \text{ kJ mol}^{-1} = 484 \text{ kJ mol}^{-1}$   
 $\Delta_r H^\circ (1) = +13 \text{ kJ mol}^{-1}$  2  
*Plus sign not required. Allow error carried forward from (i).*
- (d) Energy stored = 242 kJ 1  
*For the sequence:  $2 \times \text{reaction (1)} + \text{reaction (2)} + 2 \times \text{reaction (3)}$  there are two moles of sulfur atoms.*  
*This sequence has an overall  $\Delta_r H^\circ = 484 \text{ kJ mol}^{-1}$*   
*All this energy is 'stored' as separate hydrogen and oxygen and can be released when these are recombined.*  
*Therefore, per mol of sulfur, the energy stored is 242 kJ.*

**Question Total 12**

2. This question is about the chemistry of tungsten

(a)

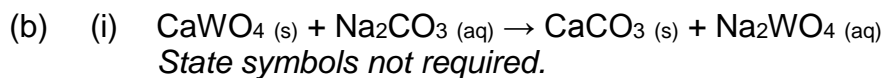


1/2

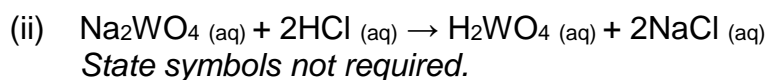
Bond Angle: 109.5°

1/2

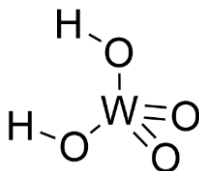
*Must have attempted to draw a tetrahedral shape to be given credit, i.e. do not credit square planar structures.*



1

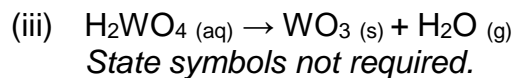


1

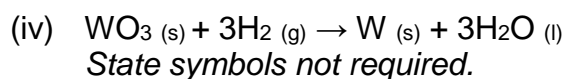


1

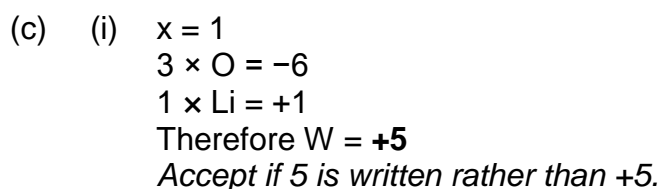
*Correct 3D tetrahedral structure not required as long as connectivity and bonding are correct.*



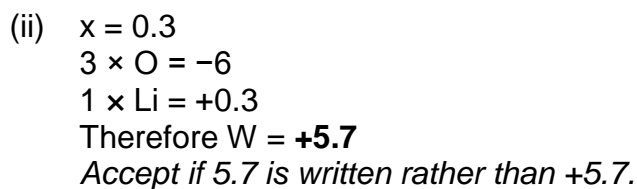
1



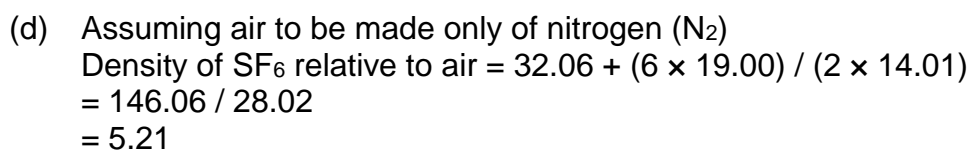
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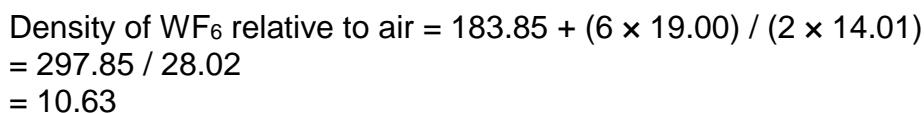
1



1



1



1

- (e)  $pV = nRT$   
 1 mol of gas occupies:  $V/n = RT/p$   
 $= (8.314 \text{ J K mol}^{-1} \times 298 \text{ K}) / 100000 \text{ N m}^{-2}$   
 $= 0.02476 \text{ m}^3 \text{ mol}^{-1}$   
 $= 2.476 \times 10^4 \text{ cm}^3 \text{ mol}^{-1}$   
 Density of  $\text{WF}_6 = 297.85 \text{ g mol}^{-1} / 2.476 \times 10^4 \text{ cm}^3 \text{ mol}^{-1}$   
 $= 0.0120 \text{ g cm}^{-3}$  2  
*Give credit if they use  $24 \text{ dm}^3$  for 1 mol of gas as a known value at STP.*
- (f)  $\text{WF}_6 (\text{g}) + 4\text{H}_2\text{O} (\text{l}) \rightarrow \text{H}_2\text{WO}_4 (\text{aq}) + 6\text{HF} (\text{aq})$  1  
 $\text{WF}_6 (\text{g}) + 3\text{H}_2\text{O} (\text{l}) \rightarrow \text{WO}_3 (\text{s}) + 6\text{HF} (\text{aq})$   
*Accept either. State symbols not required.*
- (g) (i) Positive 1  
*This is because there are more moles of gas on the right than the left.*
- (ii)  $\Delta_r H^\circ = \Delta_f H^\circ (\text{H}_2\text{SO}_4 (\text{g})) + 6 \times \Delta_f H^\circ (\text{HF} (\text{g})) - \Delta_f H^\circ (\text{SF}_6) - 4 \times \Delta_f H^\circ (\text{H}_2\text{O})$  1  
 $= -735 + (6 \times -273) - (-1210 + 4 \times -242) \text{ kJ mol}^{-1}$   
 $= -195 \text{ kJ mol}^{-1}$  1  
*1 mark for correct expression if numerical calculation is done incorrectly. Correct answer scores full marks.*
- (iii) **B**  $\text{SF}_6$  is kinetically stable but thermodynamically unstable 1
- (h) Tungsten = 1 atom inside unit cell + 4 × atoms on face + 8 × atoms on corner  
 $= 1 + (4 \times \frac{1}{2}) + (8 \times \frac{1}{8}) = \mathbf{4 \text{ atoms}}$  1
- Calcium = 6 × atoms on face + 4 × atoms on edge  
 $= (6 \times \frac{1}{2}) + (4 \times \frac{1}{4}) = \mathbf{4 \text{ atoms}}$  1
- Oxygen = 16 × atoms inside unit cell  
 $= \mathbf{16 \text{ atoms}}$  1
- (i) Volume of Unit Cell =  $0.524 \text{ nm} \times 0.524 \text{ nm} \times 1.137 \text{ nm}$   
 $= 3.122 \times 10^{-28} \text{ m}^3 = 3.122 \times 10^{-22} \text{ cm}^3$
- Molar mass of  $\text{CaWO}_4 = (40.08 + 183.85 + 4 \times 16.00) \text{ g mol}^{-1}$   
 $= 287.93 \text{ g mol}^{-1}$
- Mass of one formula unit =  $287.93 \text{ g mol}^{-1} / 6.02 \times 10^{23} \text{ mol}^{-1}$   
 $= 4.783 \times 10^{-22} \text{ g}$  1
- Mass of one unit cell  
 $= 4 \times 4.783 \times 10^{-22} \text{ g} = 1.913 \times 10^{-21} \text{ g}$  1
- Density =  $1.913 \times 10^{-21} \text{ g} / 3.122 \times 10^{-22} \text{ cm}^3$   
 $= 6.13 \text{ g cm}^{-3}$  1
- 1 mark for calculation of mass of formula unit, 1 mark for four formula units per unit cell and 1 mark for answer. Correct answer scores full marks.*

**Question Total    23**

### 3. This question is about Double Bond Equivalents, DBE

- (a) (i)  $C_nH_{2n}$   $\frac{1}{2}$   
 (ii)  $C_nH_{2n-2}$   $\frac{1}{2}$   
 (iii)  $C_nH_{2n-2}$   $\frac{1}{2}$   
 (iv)  $C_nH_{2n-6}$   $\frac{1}{2}$

(b)

DBE = 3	Ring	Double Bond	Triple Bond
	3	0	0
	2	1	0
	1	2	0
	1	0	1
	0	3	0
	0	1	1

**If all correct** (in any order) **2**  
 Minus  $\frac{1}{2}$  mark for any missing or incorrect line down to 0

DBE = 4	Ring	Double Bond	Triple Bond
	4	0	0
	3	1	0
	2	2	0
	2	0	1
	1	3	0
	1	1	1
	0	4	0
	0	2	1
	0	0	2

**If all correct** (in any order) **3**  
 Minus  $\frac{1}{2}$  mark for any missing or incorrect line down to 0

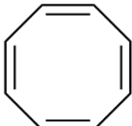

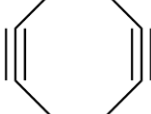
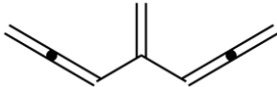
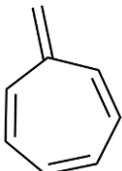
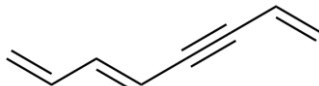
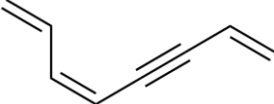
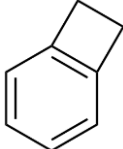
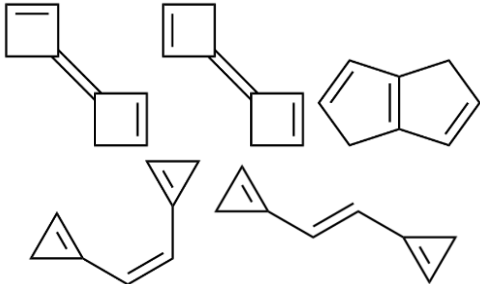
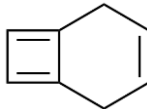
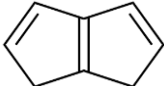
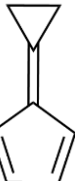
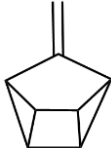
- (c) (i) 4  $\frac{1}{2}$   
 (ii) 9  $\frac{1}{2}$   
 (iii) 61  $\frac{1}{2}$   
 (iv) 4  $\frac{1}{2}$   
 (v) 4  $\frac{1}{2}$   
 (vi) 3  $\frac{1}{2}$

(d)

Spectrum	Number of atoms in each region (must add to 8)					Structural information deduced		
	Triple Bond (Alkyne)	Double Bond (Alkene)	Single Bond	Allene Central	Allene Flanking	Number of Triple Bonds	Number of Double Bonds	Number of Rings
A	0	8	0	0	0	0	4	1
B	0	0	8	0	0	0	0	5
C	4	0	4	0	0	2	0	1
D	0	2	0	2	4	0	5	0
E	0	8	0	0	0	0	4	1
F	2	6	0	0	0	1	3	0
G	0	6	2	0	0	0	3	2
H	0	6	2	0	0	0	3	2
I	0	2	6	0	0	0	1	4

$\frac{1}{2}$   
 $\frac{1}{2}$   
 $\frac{1}{2}$   
**1**  
 $\frac{1}{2}$   
 $\frac{1}{2}$   
 $\frac{1}{2}$   
 $\frac{1}{2}$   
 $\frac{1}{2}$

*Each line must be fully correct to score the mark*

<p><b>A</b></p> 	<p><b>B</b></p> 	2
<p><b>C</b></p> 	<p><b>D</b></p> 	4
<p><b>E</b></p> 	<p><b>F</b></p>  <p><i>This is the correct structure</i></p>  <p><i>This structure also fits the NMR data</i></p>	4
<p><b>G</b></p>  <p><i>This is the correct structure</i></p>	 <p><i>These structures also fit the NMR data</i></p>  <p><i>This structure is unstable as is antiaromatic but also fits the NMR data so give full credit</i></p>	2
<p><b>H</b></p>  <p><i>This is the correct structure</i></p>  <p><i>This structure also fits the NMR data</i></p>	<p><b>I</b></p> 	4

**Structures A and B are worth 1 mark each, Structures C-I are worth 2 marks each.**

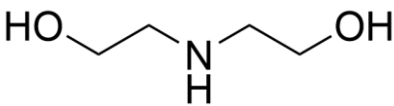
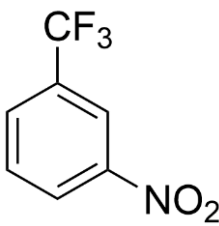
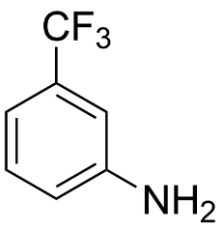
Marks are awarded for each fully correct structure in the correct place. Correct structures in an incorrect place score zero. No partial marks are awarded for a structure. No error carried forward is allowed if structure is wrong but consistent with the student's answer in the previous table. Where more there is more than one possibility only one structure needs to be drawn. There may be other possibilities which can be given full credit but only if they are **fully** consistent with all NMR data listed. Please contact the Committee if you find any alternatives.

**4 This question is about the synthesis of Addyi**

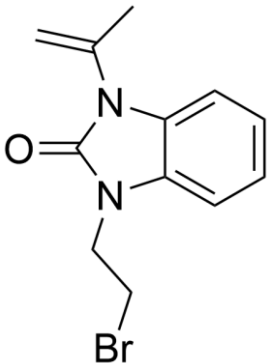
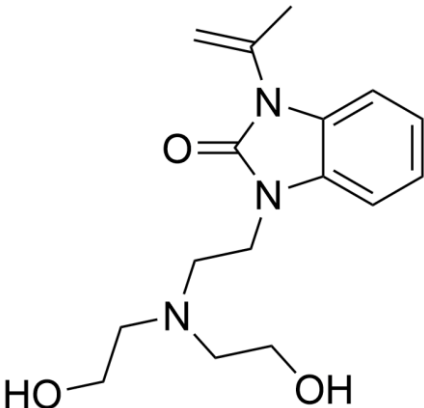
- (a) Carbon:  $45.70/12.01 = 3.805$   
 Hydrogen  $10.55/1.008 = 10.47$   
 Nitrogen  $13.32/14.01 = 0.951$   
 Oxygen  $30.43/15.99 = 1.90$   
 Simplest whole number ratio = 4:11:1:2  
 Empirical formula =  $C_4H_{11}O_2N$

*If oxygen is forgotten then can award 1 mark if calculation is done correctly.*

**2**

<p>(b) <b>Compound A</b></p>  <p><i>1 mark</i></p> <p><i>No carry forward if they propose a structure that matches their incorrect empirical formula in (a).</i></p>	<p><b>Compound B</b></p>  <p><i>1 mark</i></p> <p><i>Wrong isomer 0 marks.</i></p>	<p><b>Compound C</b></p>  <p><i>1 mark</i></p> <p><i>Allow error carried forward if wrong isomer is drawn for B and same wrong isomer is drawn for C but functional group change is correct.</i></p>
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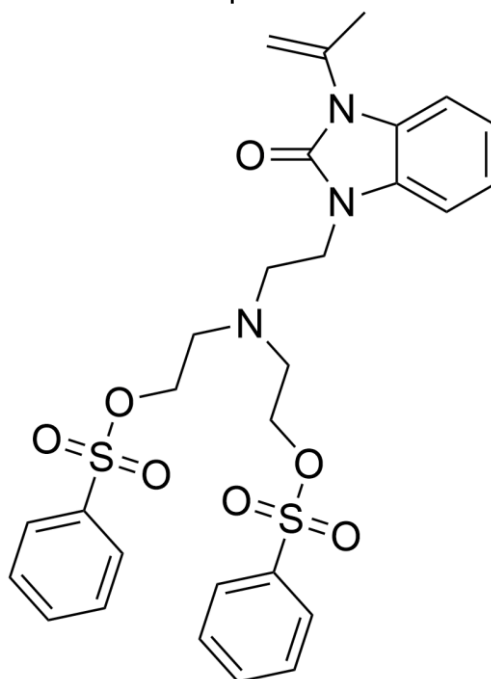
**3**

<p>(c) <b>Compound D</b></p>  <p><i>2 marks</i></p>	<p><b>Compound E</b></p>  <p><i>2 marks</i></p>
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**4**



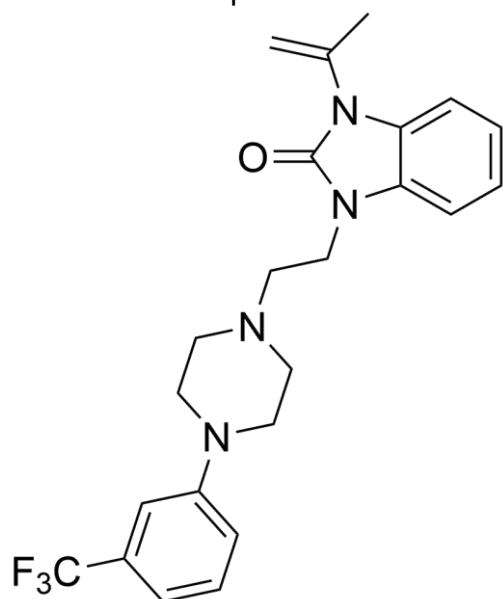
Compound **F**



2

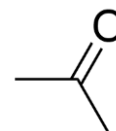
2 marks

Compound **G**



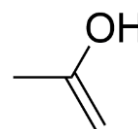
2 marks

By-Product **H**



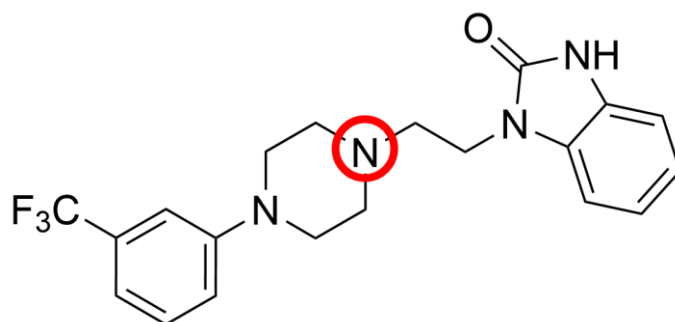
2 marks

Accept the following alternative for 1 mark.



4

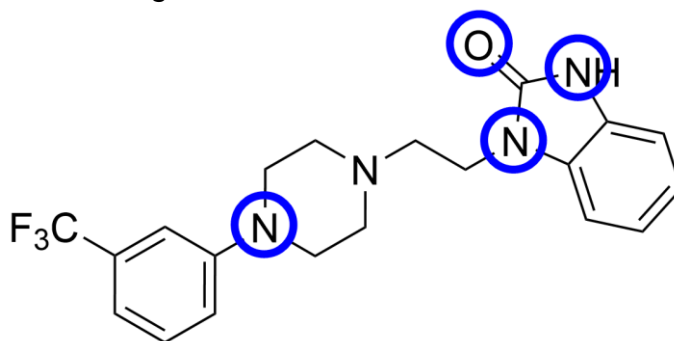
(d)



2

2 marks

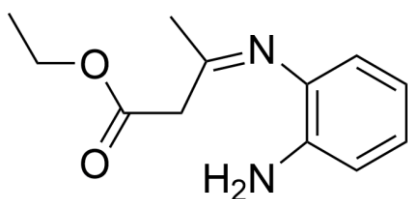
Accept any one of the following for 1 mark.



If more than one atom circled then no marks are awarded.

(e)

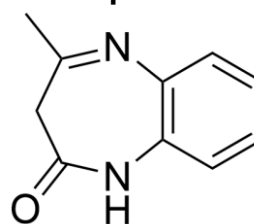
**Intermediate X**



2 marks

Give full credit to the E isomer of the imine.

**Compound Y**

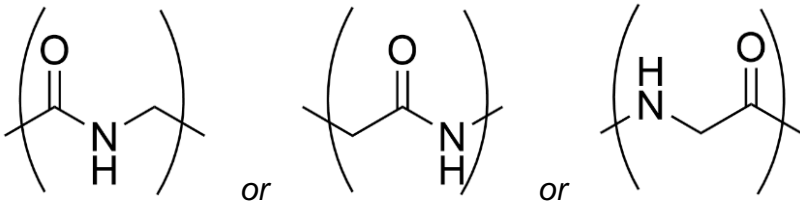


4

2 marks

**Question Total 21**

5. This question is about the radiocarbon dating of King Richard III

- (a) **C** Wash repeatedly with dilute HCl then dilute NaOH 1
- (b) (i)  1  
*Accept if they label the diagram with an 'n' to show the polymer*
- (ii)  $[\text{NHCH}_2\text{CO}]_n + 3\frac{1}{4} n\text{O}_2 \rightarrow 2n\text{CO}_2 + n\text{NO}_2 + \frac{3}{2} n\text{H}_2\text{O}$  2  
 $[\text{NHCH}_2\text{CO}]_n + 2\frac{1}{4} n\text{O}_2 \rightarrow 2n\text{CO}_2 + \frac{n}{2}\text{N}_2 + \frac{3}{2} n\text{H}_2\text{O}$   
*Accept either equation or a multiple of either.*
- (iii)  $\text{CO}_2(\text{g}) + 2\text{H}_2(\text{g}) \rightarrow \text{C}(\text{s}) + 2\text{H}_2\text{O}(\text{g})$  1  
*State symbols not required.*
- (iv) 40% of 1.0 g bone = 0.40 g polyglycine  
 The entire mass of carbon in polyglycine becomes graphite  
 Mass of graphite = % carbon in polyglycine  $\times$  0.40 g  
 $(2 \times 12.01) / (2 \times 12.01 + 3 \times 1.008 + 16.00 + 14.01) \times 0.40 \text{ g}$   
 = 0.168 g 1
- (c) Amount of C =  $0.002 \text{ g} / 12.01 \text{ g mol}^{-1} = 1.665 \times 10^{-4} \text{ mol}$   
 Number of atoms of C =  $1.665 \times 10^{-4} \text{ mol} \times 6.02 \times 10^{23} \text{ mol}^{-1} = 1.00 \times 10^{20}$   
 Number of atoms of  $^{14}\text{C}$  initially ( $N_0$ ) =  $(1.215 \times 10^{-10} / 100) \times 1.00 \times 10^{20} = 1.22 \times 10^8$  1
- Half life ( $t_{1/2}$ ) = 5568 years =  $5568 \times 365 = 2.032 \times 10^6 \text{ days}$   
 Decay Constant ( $k$ ) =  $\ln 2 / t_{1/2} = \ln 2 / 2.032 \times 10^6 \text{ days}$   
 =  $3.41 \times 10^{-7} \text{ day}^{-1}$  1
- Number of atoms of  $^{14}\text{C}$  left =  $N_0 \exp(-kt)$   
 Number of atoms of  $^{14}\text{C}$  decayed =  $N_0 - N_0 \exp(-kt)$   
 =  $1.22 \times 10^8 - 1.22 \times 10^8 \times \exp(-3.41 \times 10^{-7} \times 1)$   
 = 41.6  
 $\approx 42$  atoms 1
- (d) Decay Constant ( $k$ ) =  $\ln 2 / t_{1/2} = \ln 2 / 5568 \text{ years} = 1.245 \times 10^{-4} \text{ year}^{-1}$  1  
 $N(^{14}\text{C})/N(^{12}\text{C}) = 1.154 \times 10^{-12}$   
 $N_0(^{14}\text{C})/N_0(^{12}\text{C}) = 1.215 \times 10^{-10} / 98.93 = 1.228 \times 10^{-12}$   
 As  $N(^{12}\text{C}) = N_0(^{12}\text{C})$   
 $N(^{14}\text{C})/N_0(^{14}\text{C}) = 1.154 \times 10^{-12} / 1.228 \times 10^{-12} = 0.9397$  1
- For  $^{14}\text{C}$ :  $N = N_0 \exp(-kt)$   
 $t = - [\ln (N(^{14}\text{C})/N_0(^{14}\text{C}))] / k$   
 =  $- [\ln (0.9397)] / 1.245 \times 10^{-4} = 500 \text{ years}$  1  
 Date of Death =  $2012 - 500 = 1512$   
*Accept if any year between 2012-2016 is used as the date of the dating experiment.* 1  
*No credit if they just write 1485!*

**Question Total** 13  
**PAPER TOTAL** 100