MARK SCHEME for the May/June 2015 series

9701 CHEMISTRY

9701/21

Paper 2 (Structured Questions AS Core), maximum raw mark 60

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Page 2	Mark Scheme	Syllabus	Paper
	Cambridge International AS/A Level – May/June 2015	9701	21

Questi	on			Mark Scheme			Mark	Total
1 (a)		sub-at	tomic particle	relative mass	relative charge			
		r	neutron	1	0		[1]	
		е	electron	1/1836	-1		[1]	
			proton	1	+1		[1]	[3]
(b)	(i)		relative to 1/1	e mass of the isotope <u>s</u> / 2 the mass of an atom of (exactly) 12 (units)	an atom(s) f ¹² C / on a scale where a	an	[1] [1]	
		isotope	number v	h the same number of pr rith different mass number nucleon number		proton	[1]	[3]
	(ii)	<u>(0.89×</u>	74)+(9.37×70	$(7.63 \times 77) + (23.77 \times 70)$	78)+(49.61×80)+(8.73	×82)	[1]	
		= 79.04	l (2 d.p.) AND	Se			[1]	[2]
(c)	(i)	Те	C1					
		47.4 128	<u>52.6</u> 35.5				[1]	
		$\frac{0.370}{0.370}$	1.48 0.370					
		1	4 s	o EF = TeC <i>l</i> ₄			[1]	
			E	mpirical Formula Mass	= 270 so MF = Te	eCl₄	[1]	[3]
(c)	(ii)	Covale	nt AND simple	e/molecular			[1]	
		low me	Iting point/rea	ction with water			[1]	[2]
((iii)		+ $3H_2O \rightarrow H_2T$ C l_4 + $2H_2O \rightarrow$	eO ₃ + 4HC <i>l</i> TeO ₂ + 4HC <i>l</i>			[1]	[1]
(d)	(i)	White f	/orange flame umes/solid /green gas dis				[1] [1] [1]	[max 2]

Page 3	Mark Scheme Syllab		
	Cambridge International AS/A Level – May/June 2015 970 ⁷	1 21	
Question	Mark Scheme	Mark	Total
(ii)	NaCl giant / lattice AND ionic SiCl ₄ simple / molecular AND covalent	[1] [1]	
	For NaCl large difference in electronegativity (of sodium/Na and chlorine/ Cl/Cl_2) (indicates electron transfer/ions)	[1]	
	For SiC4 smaller difference (indicates sharing/covalency) with (weak) van der Waals'/IM forces (between molecules) ora	[4]	F 41
		[1]	[4] [20]
2 (a) (i)	Straight line drawn horizontally from same intercept	[1]	[1]
(ii)	T_1 because it shows greatest deviation/furthest from ideal	[1]	[1]
(iii)	reducing T (reduces KE of particles) so intermolecular forces of attraction become more significant	[1]	[1]
(iv)	greatest deviation is at high pressure	[1]	
	increasing pressure decreases volume so volume of particles becomes more significant ora	[1]	[2]
(b)	Mass of air= 100×0.00118 = 0.118 gMass of flask= $47.930 - 0.118$ = 47.812 gMass of Y= $47.989 - 47.812$ = 0.177 g	[1] [1]	
	$pV = nRT = \frac{m}{M_r} RT$		
	$M_r = \frac{mRT}{pV} = \frac{0.177 \times 8.31 \times 299}{1 \times 10^5 \times 100 \times 10^{-6}}$	[1]	
	= 44.0 (43.979 to 2 or more sf)	[1]	[4]
(c) (i)	strong <u>triple</u> bond	[1]	[1]
(ii)	high temperature (needed for reaction between N_2 and O_2)	[1]	[1]
(iii)	$\begin{array}{l} 2\text{NO} + 2\text{CO} \rightarrow \text{N}_2 + 2\text{CO}_2 \\ \textbf{OR} \ 2\text{NO} + \text{C} \rightarrow \text{N}_2 + \text{CO}_2 \end{array}$	[1]	[1]
(iv)	$4NO_2 + 2H_2O + O_2 \rightarrow 4HNO_3$	[1]	[1]
(v)	$NO + \frac{1}{2}O_2 \rightarrow NO_2$	[1]	
	$\begin{array}{l} NO_2 + SO_2 \rightarrow NO + SO_3 \\ \mathbf{OR} \ NO_2 + SO_2 + H_2O \rightarrow NO + H2SO_4 \end{array}$	[1]	[2]
			[15]

Page 4	Mark Scheme	Syllabus	Paper
	Cambridge International AS/A Level – May/June 2015	9701	21

Question		Mark Scheme		Total
3	(a)	Bond breaking = $C=O = 740$ C-H = 410 = 1150 kJ	[1]	
		Bond forming = $\begin{array}{c} C-C &= 350 \\ C-O &= 360 \\ O-H &= 460 \\ \end{array} = 1170 kJ$	[1]	
		Enthalpy change = $1150 - 1170 = -20 \text{ kJ mol}^{-1}$	[1]	[3]
	(b) (i)	Stereoisomerism = (molecules with the same molecular formula and) same structural formula but different spatial arrangements of atoms	[1]	
		Chiral centre = atom with four different atoms/groups attached	[1]	[2]
	(ii)	(Planar) carbonyl so (equal chance of nucleophile) attacking either side	[1]	[1]
3	(c) (i)			
		M1 = lone pair AND curly arrow from lone pair to carbonyl C M2 = partial charges on C=O AND curly arrow from bond (=) to O^{δ^-} M3 = structure of intermediate including charge M4 = lone pair AND two correct curly arrows (from lone pair to H AND from H—C to C)	[1] [1] [1] [1]	
		$M5 = CN^{-}$	[1]	[5]
	(ii)	(CN ⁻ regenerated so) catalyst	[1]	[1]
				[12]

Question Mark Scheme Mark Total 4 (a) $A = 0$ OH $A = 0OH$ $A = 0A = 0$ $A = 0A = 0$ $A = 0A = 0$ $A = 0$ <	Page 5	Mark SchemeSyllabuCambridge International AS/A Level – May/June 20159701	s Pap 21	
(b) (i) but-1-ene/1-butene but-2-ene/2-butene but	Question	Mark Scheme		
$\begin{array}{c c} & & & & & & & & & & & & & & & & & & &$	4 (a)	OH chain OH	[1]	
$\begin{array}{c c} & & & & & & & & & & & & & & & & & & &$			[1]	
$\begin{array}{ c c c c c } \hline & & & & & & & & & & & & & & & & & & $		C= OH chain OH	[1] [1] [1]	
(b) (i)but-1-ene/1-butene but-2-ene/2-butene[1] (1][2](ii)but-2-ene AND two different groups on each carbon (of C=C) double bond means no free rotation[1] [1][2](iii) $H_{-} + H_{-} + H_{$		chain OR position		
but-2-ene/2-butene[1][2](ii)but-2-ene AND two different groups on each carbon (of C=C) double bond means no free rotation[1][2](iii)HHHH[1][2](iii)HHHH[1][1](iii)HHHH[1][1](iii)HHHH[1][1](iii)HHHH[1][1](iii)HHHH[1][1](iii)IIII(iii)HHHHII(iii)HHHHII(iii)IIIII(iii)IIIII(iii)IIIII(iii)IIIII(iiii)II<		OH Chain D=		[7]
double bond means no free rotation[1][2](iii) H_{H}	(b) (i)		[1] [1]	[2]
$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	(ii)	but-2-ene AND two different groups on each carbon (of C=C) double bond means no free rotation	[1] [1]	[2]
and (either way round) [2]	(iii)		[1+1]	
[13]		and (either way round)		