

CONFIDENTIAL

January 2006

ADVANCED GCE UNIT						
MARK SCHEME						
VERSION: POST-STANDARDISATION						
MAXIMUM MARK: 90						
Syllabus / Component: 2848/01						
Chemistry: Chemistry of Natural Resources						
Paper Set Date: 11/01/06						

SUBJECT OFFICER: Steven Evans

CHECKED BY	APPROVED (Tick or initials or signature)	DATE

ADVICE TO EXAMINERS ON THE ANNOTATION OF SCRIPTS

- 1. Please ensure that you use the **final** version of the Mark Scheme. You are advised to destroy all draft versions.
- 2. Please mark all post-standardisation scripts in red ink. A tick () should be used for each answer judged worthy of a mark. Ticks should be placed as close as possible to the point in the answer where the mark has been awarded. The number of ticks should be the same as the number of marks awarded. If two (or more) responses are required for one mark, use only one tick. Half marks (½) should never be used.
- 3. The following annotations may be used when marking. <u>No comments should be written on</u> scripts unless they relate directly to the mark scheme. Remember that scripts may be returned to Centres.
 - x = incorrect response (errors may also be underlined)
 - ^ = omission mark
 - bod = benefit of the doubt (where professional judgement has been used)
 - ecf = error carried forward (in consequential marking)
 - con = contradiction (in cases where candidates contradict themselves in the same response)
 - sf = error in the number of significant figures
- 4. The marks awarded for each <u>part</u> question should be indicated in the margin provided on the right hand side of the page. The mark <u>total</u> for each question should be ringed at the end of the question, on the right hand side. These totals should be added up to give the final total on the front of the paper.
- 5. In cases where candidates are required to give a specific number of answers, (e.g. 'give three reasons'), mark the first answer(s) given up to the total number required. Strike through the remainder. In specific cases where this rule cannot be applied, the exact procedure to be used is given in the mark scheme.
- 6. Correct answers to calculations should gain full credit even if no working is shown, unless otherwise indicated in the mark scheme. (An instruction on the paper to 'Show your working' is to help candidates, who may then gain partial credit even if their final answer is not correct.)
- 7. Strike through all blank spaces and/or pages in order to give a clear indication that the whole of the script has been considered.
- 8. An element of professional judgement is required in the marking of any written paper, and candidates may not use the exact words that appear in the mark scheme. If the science is correct <u>and</u> answers the question, then the mark(s) should normally be credited. If you are in doubt about the validity of any answer, contact your Team Leader/Principal Examiner for guidance.

Abbreviations, annotations and conventions used in the Mark Scheme	/ point ; NOT () ecf AW ora	 alternative and acceptable answers for the same marking separates marking points answers which are not worthy of credit words which are not essential to gain credit (underlining) key words which <u>must</u> be used to gain credit error carried forward alternative wording or reverse argument
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Mark Sc Page 1 of 4	ChemeUnit CodeSessionYearVersion42848Jan2006Pre-State		n Ind			
Question	Expected An	swers				Marks
1 a i	air/ oxygen/O ₂ (plus nothing else) (1); (blister) copper/Cu <i>NOT pure copper</i> (IGNORE +sulphur dioxide/SO ₂ /slag <i>only</i>) (1)					2
1 a ii	liquid/(I)/molte	n				1
1 a iii	1% = 1 x 1000	$0000/100 = 10^4/10000$	(1)			1
1 a iv	purification/re	fining/electrolysis/mak	king brass NO	T smelting		1
1 b i	+1 (1); 0 (1);	0 (1); +4 (1) ALLOV	V 1+ and 4+ or	1 and 4 for (1)		4
1 b ii	redox/ oxidation	on/ reduction				1
1 b iii	copper(I) sulp	hide <i>ignore gaps</i> ecf	from oxidation	state in (i)		1
1 c i	11 electrons (1) 4s ¹ 3d ¹⁰ or reverse	d (1);			2
1 c ii	d/allow D IGI	NORE transition meta	ls			1
1 d i	methyl orange/ phenolphthalein allow small spelling errors and other suitable acid-base indicators. NOT Universal indicator or litmus					1
1 d ii	$21.2 \times 0.00100/1000 (1) = 2.12 \times 10^{-5} ecf$ if one error (1)					2
1 d iii	same answer	as (ii)				1
1 d iv	Answer to d(ii Do not accept	i) x 40 [2.12 x 10 ⁻⁵ x t rounding to 1sf.	1000/25 = 8.48	x 10 ^{−4} mol dm [−]	³] (1)	1
1 e	sulphur dioxid	e/SO ₂ ; plus three fro	om			4
	dissolves in ra damages tree reacts with ox converts it to s	ain/causes acid rain; s/plants/ lakes/ fish/bu ygen and water sulphuric acid	uildings/human	health/ leaches	aluminium;	1
	IGNORE gree	enhouse gases and pl	notochemical sn	nog		
	QWC: 2 sente ALLOW "Sulp Underline firs	ences, spelling, punctu hur dioxide" alone at a st two errors.	uation and gram the start.	mar correct (1	error) see note	
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2 a	(drain)pipes/ window frames (AW)/ doors/ roofing Must be part of fabric of house.						
2 b i	(1 -)chloroethene	correct spelling required				1	
2 b ii 2 b iii	electrophilic (1); a	ddition (1) extra selectior	ns are CON			2	
2 b iii 2 b iv	addition (polymeri	sation) NOT additional				1	
2 c i	permanent dipole	-(permanent) dipole no c	others			1	
2 c ii	—€H₂−CH Ω	-CH ₂ -CH-CH ₂ - β -	+, ∂ – correct (1); indication traction between correct	on of		2	
	−−−CHz−Ç+ α	-CH2-CH-CH2- α	oups(1)				
2 d i	H CIHHH 	H (0 C - C - or I - I al H CI al DCH ₃ co	r different order) (2); ne error (1) low less displayed low monomer units reve onnection to OCOCH ₃ m	rsed ust be thro O		2	
2 d ii	copolymer					1	
2 e	2 e chains can slide/move over each other (AW implying relative movement) (1) intermolecular / permanent dipole-permanent dipole (<i>allow abbreviations</i>) forces weaker (NOT fewer) (1) working over longer distance (AW implying greater separation) (1)					3	
2 f i	hydrogen (1); Ni,	hot/ Pt (room t and p ALL	OW high t) (1) second m	nark depends	on first	2	
2 f ii	ii primary (1) as OH attached to CH_2/C with OH attached to one other C/OH at end of chain/only one R group (1)					2	
2 f iii	aldehyde ALLOV	/ carbonyl				1	
2 f iv	 (potassium/sodium) dichromate/ correct formula (1); (sulphuric) acid IGNORE conc (1) heat/ raised temp (ALLOW reflux) provided dichromate mentioned(1); 				3		
2 g i	$M_{\rm r}$ vinyl chloride =	62.5, ethanol = 46 (1);				1	
2 g ii 2 g iii	moles vinyl chloride = 10/62.5 (= 0.16) moles ethanol = 1.5/46 (= 0.0326/0.033) ecf (1) i = moles vinyl chloride (0.16)				1		
2 g iv	 % = 0.0326 x 100/0.16 = 20% (ecf, eg ALLOW 21% if 0.033 moles ethanol used) (1) 2 sf (1) mark separately provided number follows from some working shown 					2	
2 h i	water (1); <i>plus on</i> catalyst with high	e <i>trom:</i> temp and press; catalys	t of sulphuric/phosphoric	acid depend	ls on first	2	
2 h ii	yield low/ more st	eps/ chlorine/hydrogen ch	nloride dangerous/ pollut	ant NOT cos	st-related	1	

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3 a i	halogenoalkan	e/bromoalkane (1) ALLOW I	naloalkane			1
3 a ii	HCC- H H H	H C — H H (1); 2-bromoprop ignore positions	ane ecf for 1-bror of dashes, comm	nopropa as and s	ne (1) paces	2
3 b i	aerosol <u>propellants</u> / blowing agents/ refrigerants (coolants)/ air conditioning <u>coolants</u> / fire extinguishers					1
3 b ii	three from not broken down broken down (<i>i</i> by <u>high energy</u> which cause ho chlorine atoms break down oz <i>two from</i> C-Br bond wea can be broken <i>or</i> contains C-I broken down/re QWC Logical a troposphere, st fission, photodi	 vn/ unreactive in troposphere AW) in stratosphere; <u>/frequency</u> uv/radiation; pmolytic fission/photodissocia /radicals; pne - catalysis implied aker (than C–CI); by lower energy/ frequency u H bonds; which cause it to be eact in troposphere/before reaction at least three words from ratosphere, uv, radiation, radissociation 	; ation v/visible/light/radi e more reactive aching stratosphe list used correctly lical, catalyse/cata	ation; ere. , alyst, ho	molytic	6
3 b iii	<i>two from</i> boiling point/volatility; flammability; toxicity/harm to health; cost <u>of</u> <u>manufacture (AW); (</u> un)reactivity/stability/ease of disposal. <i>IGNORE greenhouse gases</i> Mark all suggestions					2
3 c i	CH ₃ CH ₂ CH ₂ OH ALLOW propa	I/C ₃ H ₇ OH/C ₃ H ₈ O. n-1-ol if formula correct in equ	uation			1
3 c ii	CH ₃ CH ₂ CH ₂ OH (<i>etc)</i> (I/aq) + H (1) for equatior (1) for state sys	I (<i>etc as above)</i> (I/aq) + HBr ₂ O (I) I correct; <i>ecf for incorrect R</i> g mbols (<i>if first mark scored</i>)	(g/aq) → CH ₃ CH aroup	H₂CH₂Br		2
3 c iii	(anhydrous) so <i>hydrated)</i> /silica	dium sulphate/ other suitable gel NOT conc sulphuric, so	e salt (unless clea da lime	rly		1
3 c iv	(fractional) dist	illation				1
3 c v	Br^{-} intermediate c	−H complete (<i>ignore</i> +) ALLOW nd to O ⁽⁺⁾ (1)	+ on C (1); bromic	de attack	< (1);	3

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4 a	alternative hydrocarbons (eg diesel)/ oxygenates/ lean burn engines/ more complete combustion (AW)/ hybrid engines/ fuel injection/oxygen sensors /reduced drag/ lighter cars (AW)					1
4 b	wind power/ wave power/ tidal power/ biomass/ nuclear/ geothermal/ hydroelectric/ solar power/cells/panels					1
4 c i	Sun (1); uv/visible / high frequency/ high energy end of spectrum/ between visible and X-ray (1)					2
4 c ii	makes bonds/molecules vibrate (more) (1); turned into kinetic energy/ move around faster which increases temperature (1)					2
4d i	Increased CO ₂ levels in troposphere/CO ₂ (g) moves <u>equilibrium</u> (position) in equation 4.1 to right* (1) Increased CO ₂ (aq) moves <u>equilibrium</u> (position) of equation 4.2 to right* (1) * or <u>equilibrium</u> producing identified products "equilibrium moves to right" scores 1 of first 2 HCO ₂ ⁻ increases (1)					3
4 d ii	rate of forward reaction = rate of back reaction (1); <i>and one from:</i> <u>concentrations</u> of reactants and products remain constant ; closed system					2
4 d iii	system not clos	ed/ CO ₂ (g) moves away fro	m surface/CO ₂ (ad	q) ionises	(AW)	1
4 e	SiO ₂ giant cova covalent bonds CO ₂ molecular <u>weak</u> intermolec <i>ALLOW permai</i> separate molec	lent/ network solid/ lattice/ v (1) <i>IGNORE "intermolecula</i> (AW) (1) cular forces (<i>can be named</i> nent dipole - permanent dip ules/ bonds in SiO ₂ are stro	vhole structure he r" and can be abbr ole)/less energy r nger (1)	eld togethe eviated needed to	er by	3
						15