



EXAM PAPERS PRACTICE

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Practice questions created by actual examiners and assessment experts

Detailed mark scheme

Suitable for all boards

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2002

XVIII

1583

Time allowed
195 Minutes

Score

/163

Percentage

%

CHEMISTRY

**OCR
AS & A LEVEL**

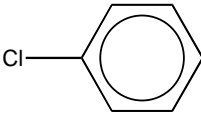
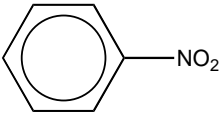
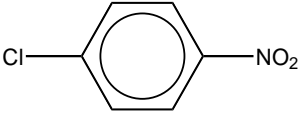
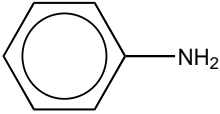
Mark Scheme

**Module 6: Organic chemistry
and analysis**

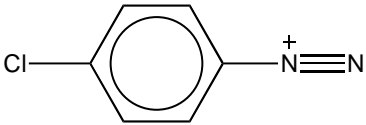
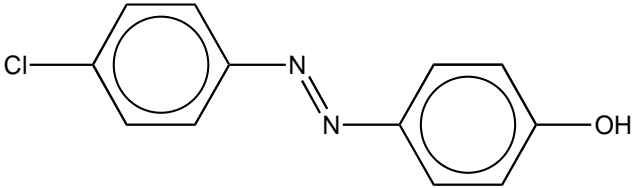
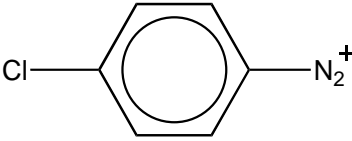
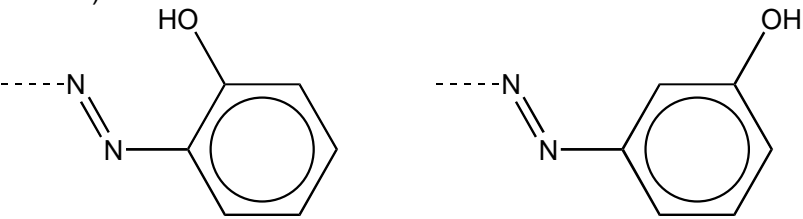
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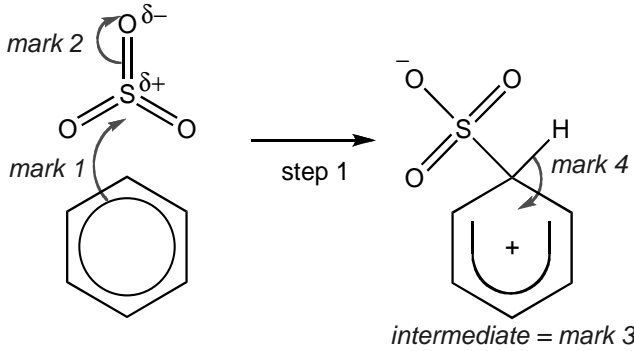
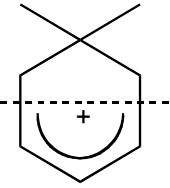
EXAM PAPERS PRACTICE

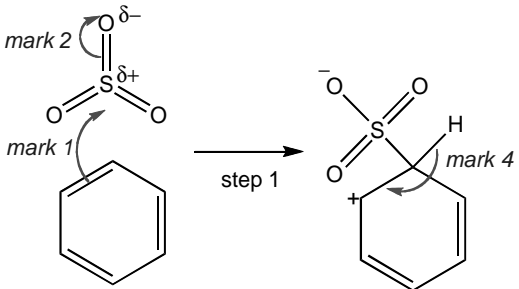
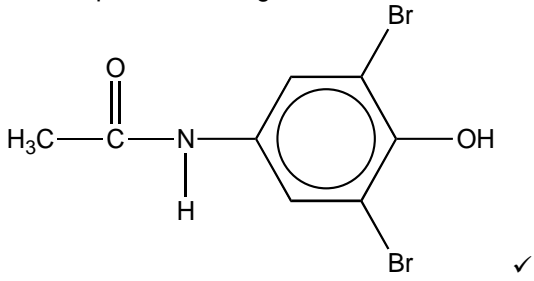
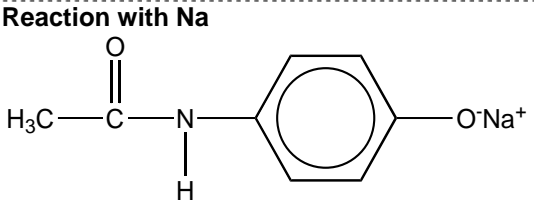
Question	er	Mark	Guidance
1 (a) (i)	<p>Response requires three stages</p> <ul style="list-style-type: none">• chlorination• nitration• reduction <p>Reduction must be a later stage than nitration</p> <p>Mark according to which sequence chosen.</p> <p>Stage 1 organic product:</p> <p> OR  ✓</p> <p>chemicals: ✓ Cl₂ AND AlCl₃ OR HNO₃ AND H₂SO₄</p> <p>Stage 2 organic product:</p> <p> OR  ✓</p> <p>chemicals: HNO₃ AND H₂SO₄ OR Sn AND HCl ✓</p> <p>Stage 3 chemicals: Cl₂ AND AlCl₃ OR Sn AND HCl ✓</p>	5	<p>Acceptable sequence of stages are:</p> <ul style="list-style-type: none">• nitration, reduction, chlorination• nitration, chlorination, reduction,• chlorination, nitration, reduction <p>For organic products, ALLOW C₆H₅NO₂ OR C₆H₅Cl OR C₆H₅NH₂ ALLOW NO₂⁻ AND NH₂⁻ DO NOT ALLOW ClC₆H₄NO₂ (formula ambiguous) DO NOT ALLOW molecular formulae IGNORE any additional structures shown eg 2- (<i>ortho</i>) and 3- (<i>meta</i>) substituted isomers</p> <p>In chemicals boxes, IGNORE temperatures IGNORE 'catalyst'</p> <p>For chlorination chemicals, ALLOW Cl₂ AND FeCl₃ OR Cl₂ AND Fe OR Cl₂ AND halogen carrier</p> <p>For nitration chemicals, 'concentrated' not required for HNO₃ OR H₂SO₄ BUT ... DO NOT ALLOW 'dilute'</p> <p>For reduction chemicals, 'concentrated' HCl not required but DO NOT ALLOW 'dilute'</p> <p>For Sn/HCl ALLOW addition of NaOH also IF it is clear that it is a second step BUT DO NOT ALLOW Sn AND HCl AND NaOH</p> <p>IGNORE catalyst</p>



Question	Answer	Mark	Guidance
(a) (i)	<p>diazonium ion</p>  <p>✓</p> <p>-N≡N group MUST be displayed</p> <p>azo dye</p>  <p>✓</p> <p>-N=N- group MUST be displayed</p>	2	<p>ALLOW '+' sign up to halfway along triple bond from left-hand N</p> <p>IGNORE presence of Cl⁻</p> <p>DO NOT ALLOW Cl⁻ substituent on benzene ring</p> <p>DO NOT ALLOW:</p>  <p>In azo dye, ALLOW as alternative to phenol OH group: O⁻ OR O⁻Na⁺ OR ONa</p> <p>ALLOW phenol part substituted at any carbon (ie 2,3 or 4 position for -OH) i.e.</p>  <p>IGNORE geometry/shape, i.e. ALLOW -N=N-</p> <p>Mark independently DO NOT ALLOW if Cl⁻ is missing from benzene ring in EITHER structure</p>



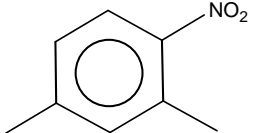
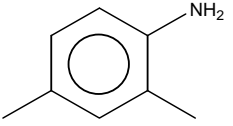
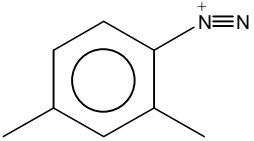
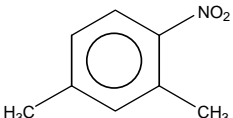
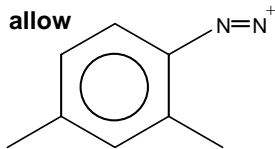
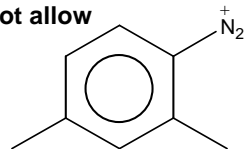
Question	er	Mark	Guidance
(b)	 <p>mark 2</p> <p>mark 1</p> <p>step 1</p> <p>intermediate = mark 3</p> <p>mark 4</p>	4	<p>ANNOTATIONS MUST BE USED</p> <p>mark 1 – curly arrow from π-delocalised ring in benzene to $S^{\delta+}$ in SO_3 ✓</p> <p>ALLOW curly arrow from the ring OR from within the ring</p> <p>mark 2 – curly arrow from one $S=O$ double bond to the O (to produce a $S-O^-$) ✓</p> <p>ALLOW curly arrow to any O in SO_3</p> <p>mark 3 – intermediate showing delocalisation over 5 carbons ✓</p> <p>Intermediate must have correct SO_3^- structure FULLY displayed</p> <p>DO NOT ALLOW intermediate with broken ring less than halfway up in correct orientation:</p>  <p>mark 4 – curly arrow from C–H bond reforming π- delocalised ring in benzene ✓</p> <p>Stand alone mark</p> <p>IGNORE responses after STEP 2</p>

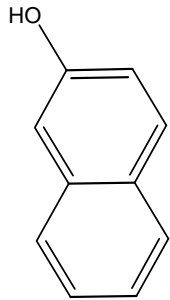
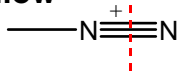
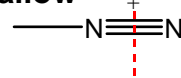
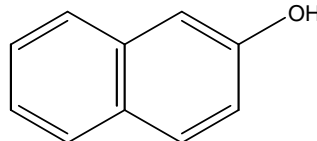
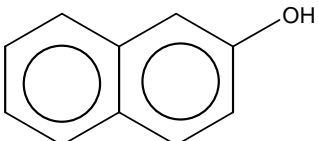
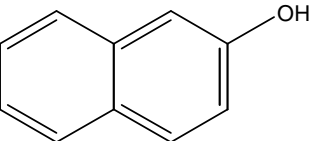
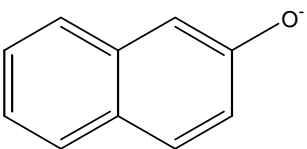
Question	er	Mark	Guidance
			<p>ALLOW Kekulé mechanism</p>  <p>ALLOW double bonds shown in other Kekulé arrangement</p>
(c)	<p>(i) Various possibilities, eg:</p>  <p>Reaction with Na</p> 	2	<p>ALLOW 1, 2, 3 or 4 Br atoms substituted on phenol ring at carbon atoms 2, 3, 5 or 6 BUT -OH must be in correct position shown DO NOT ALLOW O⁻ or ONa</p> <p>ALLOW for side chain: CH₃CONH but aromatic part of structure must be shown</p> <p>IGNORE any additional inorganic products in boxes (even if incorrect)</p> <hr/> <p>ALLOW ONa OR O⁻ as alternative to O⁻Na⁺ DO NOT ALLOW O-Na OR O⁻Na (i.e. Na without charge)</p> <p>-ONa must be in correct position shown</p> <p>ALLOW for side chain: CH₃CONH but aromatic part of structure must be shown</p> <p>IGNORE any additional inorganic products in boxes (even if incorrect)</p>

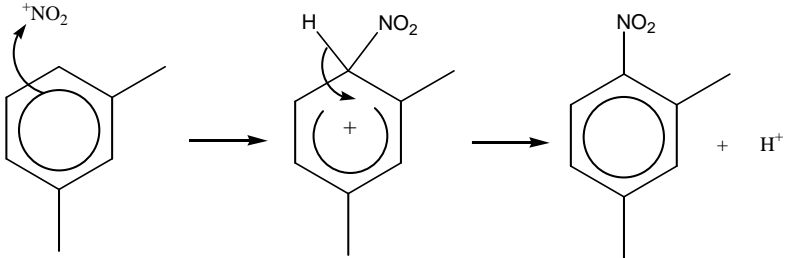
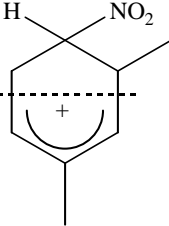


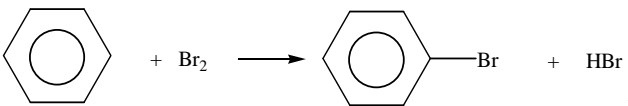
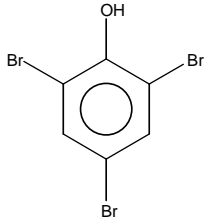
Question		er	Mark	Guidance
(c)	(ii)	<p>Hydrolysis with NaOH(aq)</p> <p>$\text{H}_3\text{C}-\overset{\text{O}}{\parallel}{\text{C}}-\text{O}^-\text{Na}^+$ ✓</p> <p>$\text{H}_2\text{N}-\text{C}_6\text{H}_4-\text{O}^-\text{Na}^+$ ✓</p> <p>Mark independently</p>	2	<p>On BOTH structures, ALLOW ONa OR O^- as alternative to O^-Na^+ DO NOT ALLOW $\text{O}-\text{Na}$ OR O^-Na (i.e. Na without charge)</p> <p>$-\text{ONa}$ must be in correct position shown on 2nd structure</p> <p>ALLOW CH_3COONa/ $\text{CH}_3\text{CO}_2\text{Na}$ OR CH_3COO^-/ CH_3CO_2^-</p> <p>ALLOW one mark for carboxylic acid AND phenol, rather than sodium salts:</p> <p>$\text{H}_3\text{C}-\overset{\text{O}}{\parallel}{\text{C}}-\text{OH}$ $\text{H}_2\text{N}-\text{C}_6\text{H}_4-\text{OH}$</p> <p>ALLOW NH_2-, CH_3-</p> <p>IGNORE any additional inorganic products in boxes (even if incorrect)</p>
Total			15	



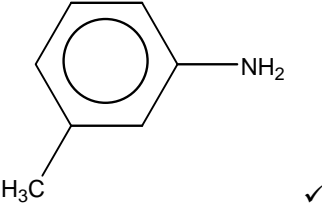
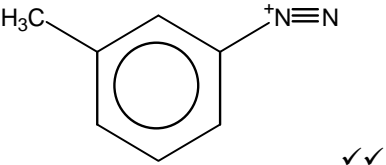
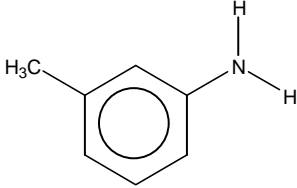
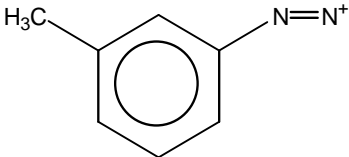
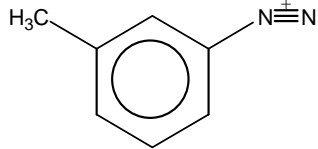
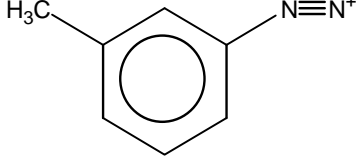
Question		Expected Answers	Marks	Additional Guidance
2	a	<p>Bond length intermediate between/different from (short) C=C and (long) C-C ✓</p> <p>ΔH hydrogenation less exothermic than expected (when compared to ΔH hydrogenation for cyclohexene) ✓</p> <p>Only reacts with Br₂ at high temp or in presence of a halogen carrier / resistant to electrophilic attack ✓</p> <p>Please annotate, use ticks to show where marks are awarded</p>	3	<p>ALLOW all carbon-carbon bonds the same length</p> <p>ALLOW ΔH hydrogenation less (negative) than expected</p> <p>ALLOW ΔH hydrogenation different from that expected</p> <p>DO NOT ALLOW ΔH halogenation/hydration</p> <p>ALLOW doesn't decolourise/react with/polarise Br₂</p> <p>ALLOW doesn't undergo addition reactions (with Br₂)</p>
	b	<p>i</p> <p>compound A</p>  <p>if NO₂ in wrong position penalise here and ECF for rest of b(i) and b(ii)</p> <p>✓</p> <p>compound B</p>  <p>✓</p> <p>compound C</p>  <p>✓</p>	4	<p>ALLOW any 4-nitro-1,3-dimethylbenzene drawn in any orientation</p> <p>ALLOW</p>  <p>drawn in any orientation</p> <p>ALLOW any 4-amino-1,3-dimethylbenzene drawn in any orientation</p> <p>ECF amine of incorrect compound A (e.g. position of NO₂ or lack of methyl sticks/groups)</p> <p>ALLOW diazonium chloride salt of 1,3-dimethylbenzene</p> <p>ECF diazonium salt/compound of incorrect compound B</p> <p>IGNORE Cl⁻ ion</p> <p>allow</p>  <p>not allow</p> 

Question	Expected Answers	Marks	Additional Guidance
	<p>compound D</p>  <p style="text-align: right;">✓</p>		<p>ALLOW if + charge is floating between the two Ns only if it is closer to the correct N</p> <p>allow  not allow </p> <p>ALLOW any of</p>     <p>ALLOW O⁻ in place of OH</p>

Question	Expected Answers	Marks	Additional Guidance
ii If NO ₂ is in correct position do not penalise even if compound A in b(i) is not in correct position	<p>mark 1 $\text{HNO}_3 + 2\text{H}_2\text{SO}_4 \rightarrow \text{H}_3\text{O}^+ + 2\text{HSO}_4^- + \text{NO}_2^+$ ✓</p> <p>mark 2 – curly arrow from π ring to $^+\text{NO}_2$ ✓</p> <p>mark 3 – intermediate with π ring broken in the correct place ✓</p> <p>mark 4 – curly arrow from C–H bond back to reform π ring AND correct products ✓</p> <p>mark 5 - $\text{H}^+ + \text{HSO}_4^- \rightarrow \text{H}_2\text{SO}_4$ ✓</p> <div style="border: 1px solid black; padding: 5px; width: fit-content; margin: 10px auto;"> Link to compound A in part (i) – cannot score full marks [in b(i) & b(ii)] if NO₂ is not adjacent to a methyl </div> 	5	<p>Equation to show formation of NO₂⁺ ion ✓</p> <p>ALLOW $\text{HNO}_3 + \text{H}_2\text{SO}_4 \rightarrow \text{H}_2\text{O} + \text{HSO}_4^- + \text{NO}_2^+$</p> <p>$\text{HNO}_3 + \text{H}_2\text{SO}_4 \rightarrow \text{HSO}_4^- + \text{H}_2\text{NO}_3^+ \rightarrow \text{H}_2\text{O} + \text{NO}_2^+$</p> <p>ALLOW mark 2 curly arrow must be from 1,3-dimethylbenzene to NO₂⁺ and ECF for marks 3 and 4</p> <p>DO NOT ALLOW intermediate</p>  <p>ALLOW CH₃s shown</p> <p>ALLOW $\text{H}_3\text{O}^+ + \text{HSO}_4^- \rightarrow \text{H}_2\text{O} + \text{H}_2\text{SO}_4$</p>
iii	2 ✓	1	No other correct response
Total		13	

Question	Expected Answers	Marks	Additional Guidance
3 (a)	 <p style="text-align: right;">✓</p>	1	<p>ALLOW $C_6H_6 + Br_2 \longrightarrow C_6H_5Br + HBr$</p> <p>DO NOT ALLOW multiple substitution DO NOT ALLOW Br^+</p>
(b) (i)	<p>White precipitate OR white solid OR white crystals ✓</p>  <p style="text-align: right;">✓</p>	2	<p>DO NOT ALLOW colourless DO NOT ALLOW white ppt <u>and</u> bubbles</p> <p>DO NOT ALLOW $Br_3C_6H_2OH$ OR 2,4,6-tribromophenol OR tribromophenol</p>
(ii)	1,2-Dibromocyclohexane ✓	1	<p>ALLOW 1,2dibromocyclohexane OR 1-2dibromocyclohexane OR 1,2dibromocyclohexane OR cyclo-1,2-dibromohexane DO NOT ALLOW dibromocyclohexane OR $C_6H_{10}Br_2$ OR structures</p>
(iii)	<p>MUST spell <u>delocalised/delocalized</u> or <u>localised/localized</u> correctly once in the answer to obtain all 5 marks</p> <p>benzene <u>electrons</u> or <u>π-bonds</u> are delocalised ✓</p> <p>phenol a <u>lone</u> or <u>non-bonded</u> pair of electrons on the oxygen or the OH group is (partially) delocalised into the ring ✓</p> <p>cyclohexene electrons are localised OR delocalised between two carbons ✓</p> <p>benzene has a lower electron density OR phenol has a higher electron density OR cyclohexene has a higher electron density ✓</p> <p>benzene cannot polarise or induce a dipole in Br_2 OR phenol can polarise the Br_2 OR cyclohexene can polarise Br_2 or the Br–Br bond ✓</p>	5	<p>ALLOW diagram to show overlap of all 6 p-orbitals for delocalisation DO NOT ALLOW benzene has delocalised structure or ring</p> <p>ALLOW diagram to show movement of lone pair into ring for phenol</p> <p>ALLOW diagram or description of overlap of 2 adjacent p-orbitals for bonding in cyclohexene DO NOT ALLOW cyclohexene has a C=C double bond IGNORE slip if cyclohexene is written as cyclohexane but π - bonding correctly described</p> <p>DO NOT ALLOW charge density OR electronegativity instead of electron density ALLOW $Br^{\delta+}$ OR electrophile Br^+ as alternate to polarise</p>



<p>(c)</p>   <p>ALLOW ECF ✓✓ on incorrect amine</p> <p>$\text{HNO}_2 + \text{HCl}$ and temp $< 10^\circ\text{C}$ OR $\text{NaNO}_2 + \text{HCl}$ and temp $< 10^\circ\text{C}$ ✓</p> <p>alkaline AND phenol (if temperature stated must be below 10°C) ✓</p>	5	<p>ALLOW</p>  <p>IGNORE Cl^- ion DO NOT ALLOW if ring is connected to the N triple bond in the diazonium or if diazonium has a negative charge ALLOW one mark for correct displayed diazonium if alkyl group is not shown</p> <p>ALLOW</p>  <p>for both marks</p> <p>ALLOW</p>  <p>for one mark</p> <p>ALLOW</p>  <p>for one mark</p> <p>ALLOW NaOH OR KOH & $\text{C}_6\text{H}_5\text{OH}$ OR phenoxide ion OR $\text{C}_6\text{H}_5\text{O}^-$ ALLOW reagents and conditions from the equations</p>
Total	14	