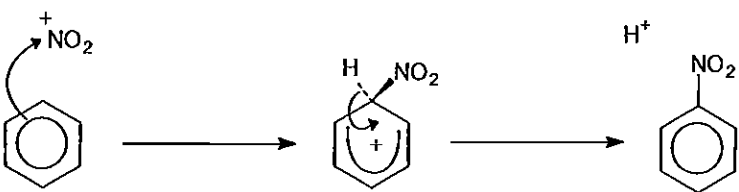


UNIT 4 API MS

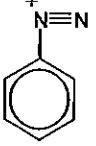
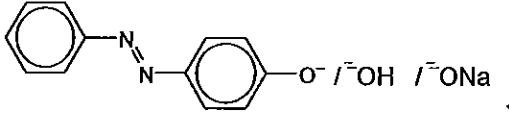
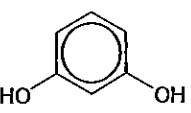
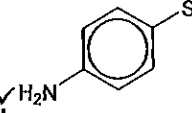
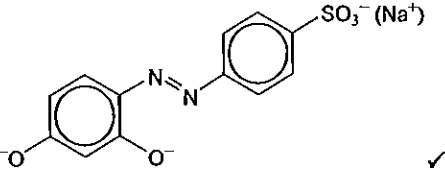
2014/01

Mark Scheme

June 2019

Qu.	Expected Answers	Marks
1 (a) (i)	$\text{HNO}_3 + \text{H}_2\text{SO}_4 \checkmark$ (both acids) conc \checkmark 50–60°C \checkmark	3
(ii)	$\text{NO}_2 \checkmark$	1
(iii)	$\text{H}_2\text{SO}_4 + \text{HNO}_3 \longrightarrow \text{NO}_2^+ + \text{H}_2\text{O} + \text{HSO}_4^- /$ $2\text{H}_2\text{SO}_4 + \text{HNO}_3 \longrightarrow \text{NO}_2^+ + \text{H}_3\text{O}^+ + 2\text{HSO}_4^- \checkmark$	1
(iv)	 <p> curly arrow from π bond to electrophile \checkmark intermediate \checkmark curly arrow from C–H bond to π bond \checkmark correct products \checkmark </p>	4
(v)	moles benzene = $\frac{3.9}{78} = 0.050 \checkmark$ actual moles of nitrobenzene formed = $\frac{4.9}{123} = 0.040 / 0.0398$ or theoretical mass nitrobenzene = $0.050 \times 123 = 6.15 \text{ (g)} \checkmark$ % yield = $\frac{\text{actual}}{\text{theoretical}} \text{ mass/moles} = 79.67\% = 80\% \checkmark$	80% without working only scores 1 mark 3
(b) (i)	$\text{AlBr}_3 / \text{Fe} / \text{FeBr}_3 \checkmark$	ALLOW AlCl_3 1
(ii)	bromine decolourised \checkmark white/cream solid/ppt. \checkmark 2,4,6-tribromophenol identified by name/structure \checkmark lone pair from O (of O–H) is delocalised into the ring (or orbital diagram to show) \checkmark increases the (π) electron density (around the ring) \checkmark Br–Br more polarised / more attracted \checkmark	ora for benzene 6
QWC	for correct use of one of the terms electrophile / electrophilic / activation	1

[Total: 20]

Qu.	Expected Answers	Marks
2 (a)	<p>step 1 $\text{HNO}_2 + \text{HCl} / \text{NaNO}_2 + \text{HCl} \checkmark$ below $10^\circ\text{C} \checkmark$</p> <p>step 2 add to phenol in alkaline conditions / NaOH (below 10°C) \checkmark</p> <div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;">  <p>\checkmark</p> </div> <div style="text-align: center;">  <p>\checkmark</p> </div> </div> <p>(b) (i) $\text{N}=\bar{\text{N}}$ circled \checkmark</p> <p>(ii) 12 carbons \checkmark 9 hydrogens \checkmark</p> <p>(iii)</p> <div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;">  <p>\checkmark</p> </div> <div style="text-align: center;">  <p>\checkmark</p> </div> </div> <p>(c)</p> <div style="text-align: center;">  <p>\checkmark</p> </div>	<p>ALLOW C_6H_5 not displayed</p> <p>ALLOW any substitution position in the dye</p> <p style="text-align: center;">5</p> <p style="text-align: center;">1</p> <p style="text-align: center;">2</p> <p>DO NOT ALLOW connection errors here</p> <p>ALLOW $-\text{SO}_3\text{H}$</p> <p style="text-align: center;">2</p> <p>ALLOW just one $\bar{\text{O}}$</p> <p style="text-align: center;">1</p>
[Total: 11]		