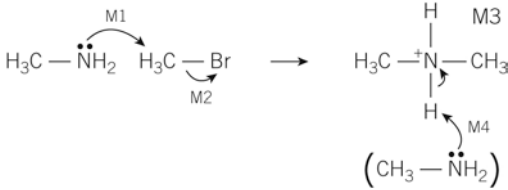
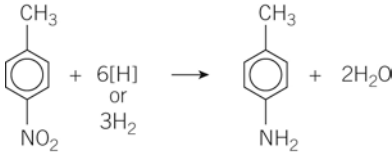
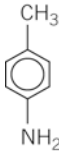
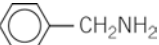
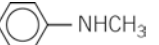
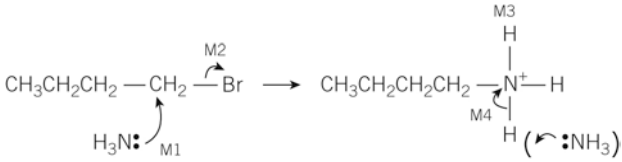
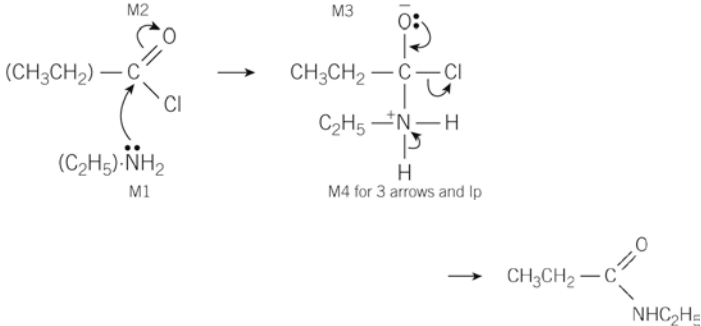


Question number	Answer	Marks	Guidance
1 (a)	Dimethylamine	1	
1 (b)	nucleophilic substitution 	5	The curly arrow must be from the lone pair. Marks 1 and 2 are for the curly arrows if shown (M1 and M2). Mark 3 is for the structure of the intermediate. Mark 4 is to show the removal of the H <sup>+</sup> . You do not need to show the reagent here. Final mark is for the name of the mechanism.
1 (c)	quaternary ammonium salt surfactant / detergent / fabric softener or conditioner/ hair conditioner	1 1	
2 (a)	Sn or Fe/HCl Sn or Fe/H <sub>2</sub> SO <sub>4</sub> OR H <sub>2</sub> /Ni can be used as the reducing agent. 	2	Do not accept dilute or conc. for these 2 acids or you cannot have HNO <sub>3</sub> at all. Do not accept NaBH <sub>4</sub> /LiAlH <sub>4</sub> or Na/C <sub>2</sub> H <sub>5</sub> OH.
2 (b)		1	NH <sub>3</sub> <sup>+</sup> is also acceptable.
2 (c)	G:  H: 	2	Do not accept the correct answers incorrectly labelled.
3 (a)	Nucleophilic substitution 	5	Mark 1, Mark 2 and Mark 4 are for the curly arrows (M1, M2 and M4). Mark 3 is for the structure of cation. Final mark is for the name of the mechanism.

3 (b)	<b>Step 1:</b> $\text{CH}_3\text{CH}_2\text{CH}_2\text{Br} + \text{KCN} \rightarrow \text{CH}_3\text{CH}_2\text{CH}_2\text{CN} + \text{KBr}$ (or $\text{CN}^-$ ) (or $\text{Br}^-$ )  <b>Step 2</b> $\text{CH}_3\text{CH}_2\text{CH}_2\text{CN} + 2\text{H}_2 \rightarrow \text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{NH}_2$	3	Do not accept HCN here.  Accept 4[H] instead of $2\text{H}_2$ .
3 (c)	There is a lone pair (on N); R group increases electron density on N / has positive inductive effect	1 1	
3 (d)	$\text{CH}_3\text{CH}_2\text{N}(\text{CH}_3)_2$	1	
4 (a)	$\text{CH}_3\text{CH}_2\text{CH}_2\text{Br}$	1	Allow Cl instead of Br.
4 (b)	$\text{CH}_3\text{CH}_2\text{CN}$	1	
4 (c)	nucleophilic substitution or from $\text{CH}_3\text{CH}_2\text{CH}_2\text{Br}$ gives less pure produce because further substitution occurs The impurity can be any one of $(\text{CH}_3\text{CH}_2\text{CH}_2)_2\text{NH}$ $(\text{CH}_3\text{CH}_2\text{CH}_2)_3\text{N}$ $(\text{CH}_3\text{CH}_2\text{CH}_2)_4\text{N}^+\text{Br}^-$	3	Also accept 'have salts including $\text{NH}_4\text{Br}$ as impurities'.
5 (a)	(nucleophilic) addition-elimination  <p>M4 for 3 arrows and lp</p>	1  4	<ul style="list-style-type: none"> <li>• Allow attack by <math>:\text{NH}_2\text{CH}_2\text{CH}_2\text{CH}_3</math></li> <li>• M2 not allowed independent of M1, but allow M1 for correct attack on C+</li> <li>• + rather than <math>\delta+</math> on C=O loses M2</li> <li>• If Cl lost with C=O breaking, max 1 for M1</li> <li>• M3 for correct structure <u>with charges</u> but lone pair on O is part of M4</li> <li>• 3 arrows in M4 can be shown in two separate steps.</li> <li>• If M3 drawn twice, mark first answer eg ignore missing + if missed off second structure</li> <li>• Only allow M4 after correct / very close M3</li> <li>• For M4, ignore <math>\text{RNH}_2</math> removing <math>\text{H}^+</math> but lose M4 for <math>\text{Cl}^-</math> removing <math>\text{H}^+</math> in mechanism,</li> <li>• but ignore HCl shown as a product.</li> </ul>
	Allow wrong amine in M1 but penalise in M3  Allow $\text{C}_3\text{H}_7$ in M3  Minus sign on $\text{NH}_3$ loses M1 (but not M4 if $\text{NH}_3$ also shown here)		

	N-propylethanamide must be this name even if wrong amine used	1	NOT N-propylethanamide
5 (b)	$\begin{array}{c} \text{H}_3\text{C}-\text{CH}-\text{CH}_3 \\   \\ \text{NH}_2 \end{array}$ Primary	1	Label and structure must both be correct for each type to score the mark. Penalize wrong number of carbons but otherwise correct, first time only Not allow ambiguous $\text{C}_3\text{H}_7\text{NH}_2$ BEWARE No mark for the original amine $\text{CH}_3\text{CH}_2\text{CH}_2\text{NH}_2$ Allow $\text{C}_2\text{H}_5$
	$\begin{array}{c} \text{H}_3\text{C}-\text{N}-\text{CH}_2\text{CH}_3 \\   \\ \text{H} \end{array}$ Secondary	1	
	$\begin{array}{c} \text{H}_3\text{C}-\text{N}-\text{CH}_3 \\   \\ \text{CH}_3 \end{array}$ Tertiary	1	
6 (a)	(nucleophilic) addition-elimination  N-ethylpropanamide	1 4 1	minus on $\text{NH}_2$ loses M1 M2 not allowed independent of M1, but allow M1 for correct attack on $\text{C}^+$ $+\text{C}=\text{O}$ loses M2 only allow M4 after correct or very close M3 lose M4 for $\text{Cl}^-$ removing $\text{H}^+$ in mechanism, but ignore HCl as a product
6 (b)	$\text{CH}_3\text{CN}$ or ethan(e)nitrile or ethanonitrile  for each step wrong or no reagent loses condition mark  Step 1 $\text{Cl}_2$ uv or above $300^\circ\text{C}$ Step 2 KCN aq and alcoholic (both needed)  Step 3 $\text{H}_2/\text{Ni}$ or $\text{LiAlH}_4$ or $\text{Na}/\text{C}_2\text{H}_5\text{OH}$	1 1 1 1 1 1	not ethanitrile but allow correct formula with ethanitrile contradiction loses mark wrong or no reagent loses condition mark allow uv light / (sun)light / uv radiation not $\text{CN}^-$ but mark on NOT HCN or KCN + acid, and this loses condition mark NOT $\text{NaBH}_4$ Sn/HCl (forms aldehyde!) ignore conditions

7 (a)	Phenylamine accepts a proton/ $H^+$	1	
7 (b)	It is an ionic salt so it dissolves in a polar solvent like water.	1	