

- M1.** (a) Cyclohexane evolves  $120 \text{ kJ mol}^{-1}$   
 $\therefore$  (expect triene to evolve)  $360 \text{ kJ mol}^{-1}$  **(1)** or  $3 \times 120$

$$360 - 208 = 152 \text{ kJ (1) NOT 150}$$

152 can score first 2

QofL: benzene lower in energy / more (stated) stable **(1)**  
 Not award if mentions energy required for bond breaking

due to delocalisation **(1)** or explained

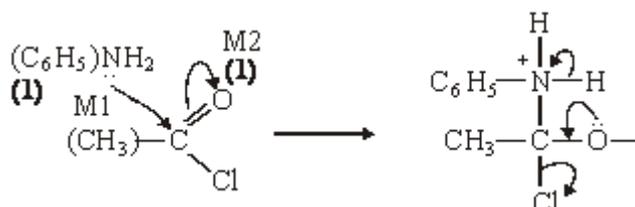
4

- (b) (i) phenylamine weaker **(1)**  
 if wrong no marks

lone pair on N (less available) **(1)**  
 delocalised into ring **(1)** or "explained"

3

- (ii) addition – elimination **(1)**



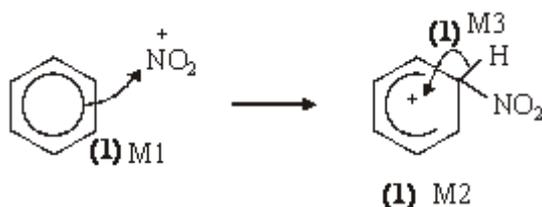
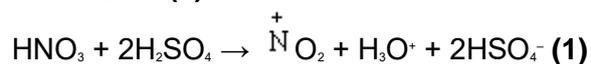
structure **(1)** M3

3 arrows **(1)** M4

N-phenyl ethanamide **(1)**

6

- (iii) conc  $\text{HNO}_3$  **(1)**  
 conc  $\text{H}_2\text{SO}_4$  **(1)**



(iv) peptide / amide (1)

NaOH (aq) (1)

*HCl conc or dil or neither*

*H<sub>2</sub>SO<sub>4</sub> dil NOT conc*

*NOT just H<sub>2</sub>O*

2

### Notes

- (a)
- 360 or  $3 \times 120$  or in words (1);
  - 152 NOT 150 (1); (152 can get first two marks)
  - **Q of L** benzene more stable but not award if  $\Delta H$  values used to say that more energy required by benzene for hydrogenation compared with the triene or if benzene is only compared with cyclohexene (1);
  - delocalisation or explained (1)

- (b) (ii) or N-phenylacetamide or acetanilide  
 mechanism: if shown as substitution can only gain M1  
 if  $\text{CH}_3\text{CO}^+$  formed can only gain M1  
 lose M4 if  $\text{Cl}^-$  removes  $\text{H}^+$   
 be lenient with structures for M1 and M2 but must be correct for M3  
 alone loses M2

- (iii) **No marks for name of mechanism in this part**  
 if conc missing can score one for both acids (or in equation)  
 allow two equations

allow  $\text{HNO}_3 + \text{H}_2\text{SO}_4 \rightarrow \text{NO}_2^+ + \text{HSO}_4^- + \text{H}_2\text{O}$

ignore side chain in mechanism even if wrong

arrow for M1 must come from inside hexagon

arrow to  $\text{NO}_2^+$  must go to N but be lenient over position of +

+ must not be too near "tetrahedral" Carbon

horseshoe from carbons 2-6 but don't be too harsh

(iv) reagent allow NaOH

HCl conc or dil or neither  
H<sub>2</sub>SO<sub>4</sub> dil or neither but not conc  
not just H<sub>2</sub>O

[21]

**M2.A**

[1]