## M1. (a) M1 CH<sub>3</sub>CH<sub>2</sub>COOH not C<sub>3</sub>H<sub>7</sub>COOH

1

M2 CH<sub>3</sub>CH<sub>2</sub>OH or C<sub>2</sub>H<sub>5</sub>OH

1

M3 CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>COOCH<sub>2</sub>CH<sub>3</sub> + H<sub>2</sub>O

allow C<sub>3</sub>H<sub>7</sub>COOC<sub>2</sub>H<sub>5</sub>

allow C₃H₁COOC₂H₅ penalise M3 for wrong products and unbalanced equation

1

**M4** H<sub>2</sub>SO<sub>4</sub> or HCl or H<sub>3</sub>PO<sub>4</sub> conc or dil or neither not HNO<sub>3</sub>

1

(b) M1 CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>OH

1

not C₄H₃OH

**M2** (CH<sub>3</sub>CO) <sub>2</sub>O

1

M3 → CH₃COOCH₂CH₂CH₂CH₃ + CH₃COOH

allow CH₃COOC₄H₅ penalise M3 for wrong products and unbalanced equation

1

(c) (nucleophilic) addition-elimination

not acylation alone

M2 not allowed indep of M1 but allow M1 for correct attack on C+

+C=O loses M2

only allow M4 after correct or v close M3 ignore CI removing H-

5

(d)

$$\begin{pmatrix}
CH_{2}OOCC_{17}H_{31} \\
CH_{2}OOCC_{17}H_{33} \\
CH_{2}OOCC_{17}H_{29}
\end{pmatrix} + 3 CH_{3}OH$$

$$CH_{2}OH$$

$$CH_{$$

ignore errors in initial triester First mark for 3CH<sub>3</sub>OH Third mark for all three esters

3

(e)

First mark for correct ester link second mark for the rest including trailing bonds

If ester link wrong, lose second mark also

2

Adv reduces landfill

saves raw materials

lower cost for recycling than making from scratch reduces CO<sub>2</sub> emissions by not being incinerated

not allow cost without qualification

ignore energy uses

1

Disad difficulty/cost of collecting/sorting/processing

product not suitable for original purpose, easily contaminated

not allow cost without qualification ignore energy uses

1

[19]

**M2.** (a) (i) **W** 3

**X** 4

1

1

1

1

1

Y 2

(ii)  $H \longrightarrow H$   $H \longrightarrow C \longrightarrow Si \longrightarrow C \longrightarrow H$   $H \longrightarrow C \longrightarrow H$ 

displayed formula shows ALL bonds

(b) (i) NO<sub>2</sub><sup>+</sup> allow + anywhere can score in equation

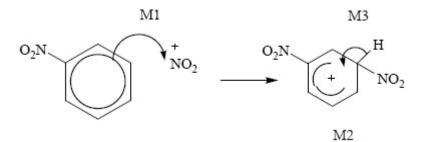
 $HNO_3 + 2H_2SO_4 \rightarrow NO_2^+ + 2HSO_4^- + H_3O^+$ 

OR

 $HNO_3 + H_2SO_4 \rightarrow NO_2^+ + HSO_4^- + H_2O$ or use two equations via  $H_2NO_3^+$ 

(ii) electrophilic substitution

Not Friedel Crafts



Allow Kekule structures

+ must be on N of 'NO<sub>2</sub> (which must be correct) both NO<sub>2</sub> must be correctly positioned and bonded to gain M2

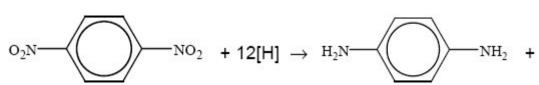
M1 arrow from circle or within it to N or to + on N horseshoe must not extend beyond C2 to C6 but can be smaller

+ not too close to C1 M3 arrow into hexagon unless Kekule allow M3 arrow independent of M2 structure ignore base removing H in M3

3

(c) (i) H<sub>2</sub>/Ni or H<sub>2</sub>/Pt or Sn/HCl or Fe/HCl (conc or dil or neither) allow dil H<sub>2</sub>SO<sub>4</sub> ignore mention of NaOH

Not NaBH<sub>4</sub> Not LiAIH<sub>4</sub> Not Na/C<sub>2</sub>H<sub>5</sub>OH not conc H<sub>2</sub>SO<sub>4</sub> or any HNO<sub>3</sub>



4H<sub>2</sub>O Or 6H<sub>2</sub>

> allow  $C_6H_4(NO_2)_2$  etc , allow  $NO_2$ —  $NH_2$  i.e. be lenient on structures, the mark is for balancing equ

1

1

ignore [], as in polymer

1st mark for correct peptide link 2nd mark for the rest correct including trailing bonds

2

(iii) **M1** Kevlar is <u>biodegradeable</u> but polyalkenes not allow Kevlar is <u>more</u> biodegradeable

1

<b>M2</b> Kevlar has <u>polar</u> bonds/is a (poly) amide/has peptide link comment on structure of Kevlar		
comment on structure of Nevial	1	
M3 can be hydrolysed/attacked by nucleophiles/acids/bases/enzymes		
	1	
M4 polyalkenes <u>non polar</u> /has <u>non-polar</u> bonds		
comment on structure of polyalkenes but not just strong bonds		
	1	[18]