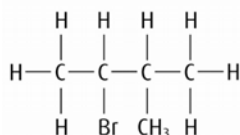
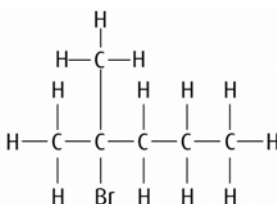


Marking scheme for Core Worksheet – Option G

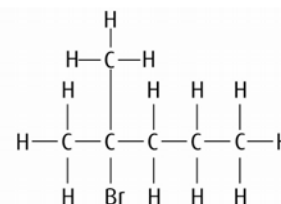
1 a



product from A



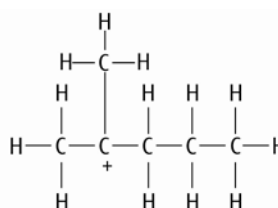
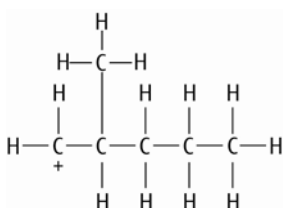
product from B



product from C

[3]

b



[2]

tertiary carbocation (on right) identified as more stable

[1]

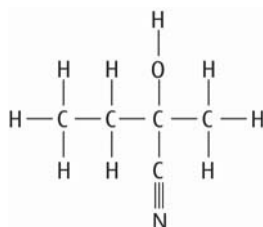
tertiary carbocation more stable than primary due to more alkyl groups around C⁺

[1]

alkyl groups electron-releasing/positive inductive effect

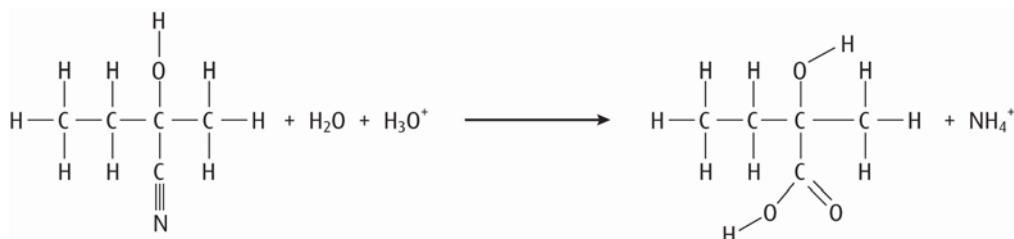
[1]

2 a



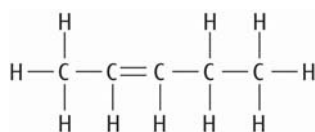
[1]

b



[2]

3 a

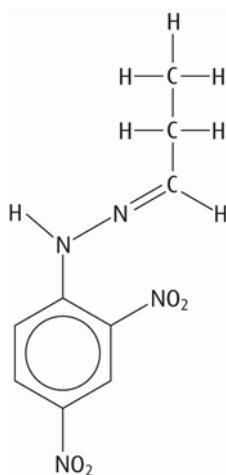


[1]

b dehydration/elimination

[1]

4



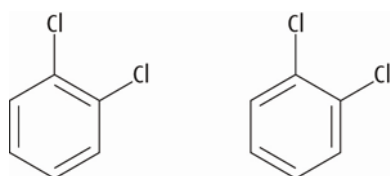
[2]

1 mark for C=N, 1 mark for rest of molecule

5

If benzene has a delocalised structure, only one 1,2-substituted isomer is possible. [1]

If benzene has the Kekulé structure, there would be two different 1,2-substituted isomers. [1]



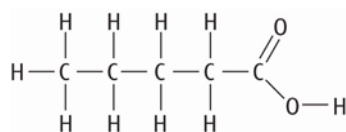
[1]

Only one 1,2-substituted isomer has been found. [1]

6

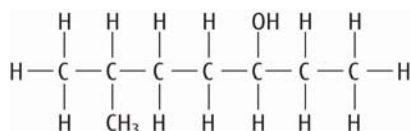
a magnesium [1]

dry ether/ethoxyethane [1]

 $\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{MgBr}$ [1]**b**

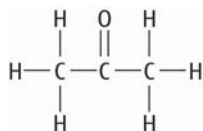
[1]

pentanoic acid [1]

c

[1]

7



[1]

 CH_3MgBr

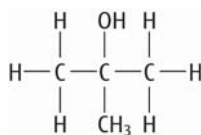
[1]

dry ether

[1]

 $\text{H}^+/\text{H}_2\text{O}$

[1]



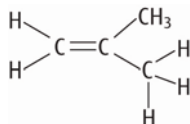
[1]

concentrated phosphoric acid

[1]

heat

[1]



[1]

8

a chloroethanoic acid is stronger

[1]

Cl is electron-withdrawing/very electronegative

[1]

electron density withdrawn from COO^- in conjugate base

[1]

 H^+ attracted less strongly/conjugate base more stable

[1]

b 4-nitrophenol is stronger

[1]

nitro group is electron-withdrawing

[1]

electron density withdrawn from O^- in conjugate base

[1]

 H^+ attracted less strongly/conjugate base more stable

[1]