## 6.3 Analysis

## 6.3.1 Chromatography and qualitative analysis

Types of chromatography				
(a) interpretation of one-way TLC chromatograms in terms of Rf values				
(b) interpretation of gas chromatograms in terms of:				
(i) retention times				
(ii) the amounts and proportions of the components in a mixture.				
Tests for organic functional groups				
(c) qualitative analysis of organic functional groups on a test-tube scale; processes and techniques needed to identify the following functional groups in an unknown compound:				
(i) alkenes by reaction with bromine				
(ii) haloalkanes by reaction with aqueous silver nitrate in ethanol				
(i) phenols by weak acidity but no reaction with CO <sub>3</sub> <sup>2-</sup>				
(ii) carbonyl compounds by reaction with 2,4- DNP				
(iii) aldehydes by reaction with Tollens' reagent				
(iv) primary and secondary alcohols and aldehydes by reaction with acidified dichromate				
(v) carboxylic acids by reaction with CO <sub>3</sub> <sup>2-</sup>				

## 6.3.2 Spectroscopy

NMR Spectroscopy			
(a) analysis of a carbon-13 NMR spectrum of an organic molecule to make predictions about:			
(i) the number of carbon environments in the molecule			
(ii) the different types of carbon environment present, from chemical			
shift values			
(iii) possible structures for the molecule			
(b) analysis of a high resolution proton NMR spectrum of an organic molecule to make predictions about:			
<ul> <li>(i) the number of proton environments in the molecule</li> <li>(ii) the different types of proton environment present, from chemical shift values</li> <li>(iii) the relative numbers of each type of proton present from relative peak areas, using integration traces or ratio numbers, when required</li> <li>(iv) the number of non-equivalent protons adjacent to a given proton from the spin- spin splitting pattern, using the n + 1 rule</li> <li>(v) possible structures for the molecule</li> </ul>			
(c) prediction of a carbon-13 or proton NMR spectrum for a given molecule			
<ul> <li>(d)</li> <li>(i) the use of tetramethylsilane, TMS, as the standard for chemical shift measurements</li> <li>(ii) the need for deuterated solvents, e.g. CDCl 3, when running an NMR spectrum</li> <li>(iii) the identification of O–H and N–H protons by proton exchange using D<sub>2</sub>O</li> </ul>			
Combined techniques			
(e) deduction of the structures of organic compounds from different analytical data including:			
(i) elemental analysis (ii) mass spectra (iii) IR spectra (iv) NMR spectra.			