

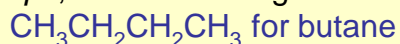
## Naming organic Chemistry A2

N Goalby  
Chemrevise.org

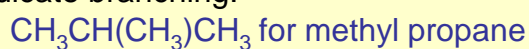
### DIFFERENT FORMULAE

- **General** represents *any member* of a homologous series  
 $C_nH_{2n+2}$  for an alkane;  
 $C_nH_{2n}$  for an alkene
- **Molecular** shows the *exact number* of atoms of each type  
 $C_4H_{10}$  for butane
- **Empirical** shows the *simplest whole number ratio* of atoms  
 $C_2H_5$  for butane

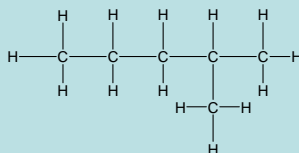
• **Condensed Structural** the minimal detail, using *conventional groups*, for an unambiguous structure.



Brackets indicate branching.

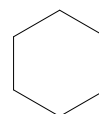


**Displayed** shows every bond and every atom

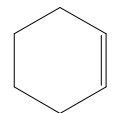


**Skeletal** used to show a simplified organic molecule. No carbon or hydrogen atoms are drawn

- A carbon atom is represented by a join of lines at an angle
- Functional groups are drawn
- The number of hydrogen atoms on each carbon atom will be how many extra bonds needed to make up four bonds



cyclohexane



cyclohexene



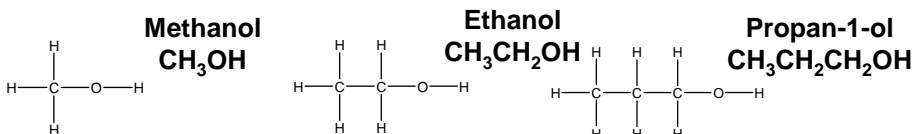
butane

## HOMOLOGOUS SERIES

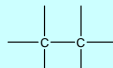
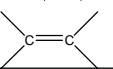
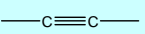
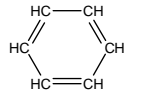

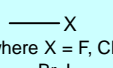
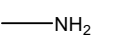
These are a series of compounds of similar structure in which each member differs from the next by a common repeating unit,  $\text{CH}_2$ . Series members are called homologues.

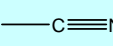
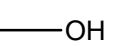
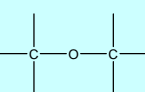
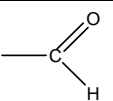
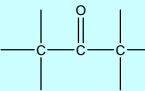
- All share the **same general formula**.
- Formula of a homologue differs from its neighbour by  $\text{CH}_2$ . (e.g.  $\text{CH}_4$ ,  $\text{C}_2\text{H}_6$ , ... etc)
- Contain the **same functional group**
- Have **similar chemical properties**.
- Show a gradual change in physical properties as molar mass increases.
- Can usually be prepared by similar methods.

E.g. the first three homologues of the alcohol series, each containing the  $-\text{OH}$  functional group



## Different types of functional groups

Group of compounds	Functional group	Naming compound (prefix-) (-suffix)	Example	
			Name	Structure
Alkane		-ane	Butane	CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>
Alkene		-ene	Propene	CH <sub>2</sub> CHCH <sub>3</sub>
Alkyne		-yne	Ethyne	CHCH
Arene			Benzene	
Haloalkane	 where X = F, Cl, Br, I	<i>Halo-</i> [Fluoro-, Chloro-, Bromo-, Iodo-]	1-bromobutane	CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> Br
Amine		-amine	Methylamine	CH <sub>3</sub> NH <sub>2</sub>

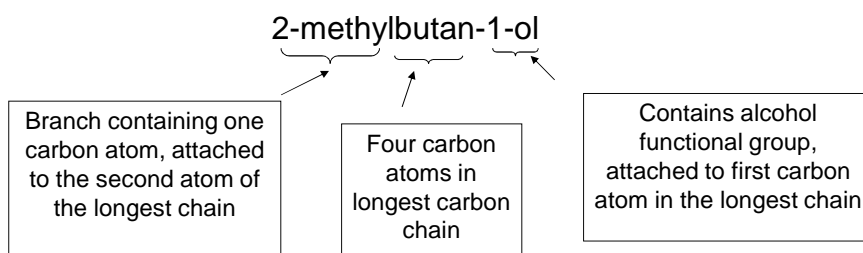
Nitrile		-nitrile	Pentanenitrile	C <sub>4</sub> H <sub>9</sub> CN
Alcohol		-ol	Propan-1-ol	CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> OH
Ether		-oxy-ane	Ethoxyethane	CH <sub>3</sub> OCH <sub>3</sub>
Aldehyde		-al	Ethanal	CH <sub>3</sub> CHO
Ketone		-one	Butanone	CH <sub>3</sub> CH <sub>2</sub> COCH <sub>3</sub>

Carboxylic acid		<i>-oic acid</i>	Ethanoic acid	CH <sub>3</sub> COOH
Carboxylate ion		<i>-oate</i>	Sodium ethanoate	CH <sub>3</sub> COO <sup>-</sup> Na <sup>+</sup>
Ester		<i>-yl -oate</i>	Ethyl ethanoate	CH <sub>3</sub> COOCH <sub>2</sub> CH <sub>3</sub>
Acyl chloride		<i>-oyl chloride</i>	Ethanoyl chloride	CH <sub>3</sub> COCl
Amide		<i>-amide</i>	Ethanamide	CH <sub>3</sub> CONH <sub>2</sub>

## NAMING ORGANIC COMPOUNDS

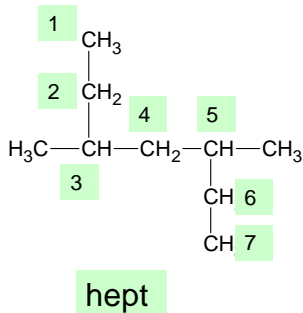
The systematic name of an organic compound tells us:

- The **number of carbon atoms** in the **longest chain** or ring.  
*This is the prefix of the name*
- The nature and position of any **functional group(s)**.  
*This is the suffix of the name*
- The position of any **side chain branches** and **number of carbon atoms in the branch**.



### Naming organic compounds: The different components

Count the number of carbon atoms in the longest chain

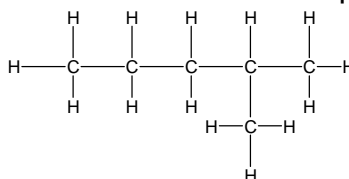


code	no of carbons
meth	1
eth	2
prop	3
but	4
pent	5
hex	6
hept	7
oct	8
non	9
dec	10

### Naming Branched Hydrocarbons

- Count the longest carbon chain and name appropriately

Eg pentane



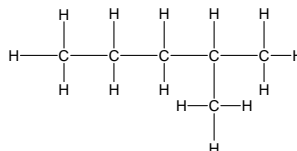
- Locate any side chains and count how many carbons they contain  
Eg  $-\text{CH}_3$  methyl or  $-\text{C}_2\text{H}_5$  ethyl

Number of carbon atoms in branch	Name of alkyl group	Structural Formula
1	Methyl-	$\text{CH}_3-$
2	Ethyl-	$\text{CH}_3\text{CH}_2-$
3	Propyl-	$\text{CH}_3\text{CH}_2\text{CH}_2-$
4	Butyl-	$\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2-$
5	Pentyl-	$\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2-$

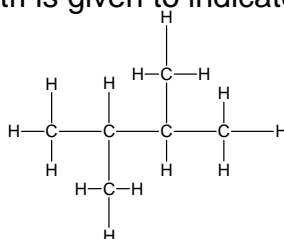
## Naming branched compounds cont.

3. Give the side chain the number the carbon is on. Eg 2-methyl. (Count from either end to give the lowest number.)

In the example above:  
2-methyl pentane

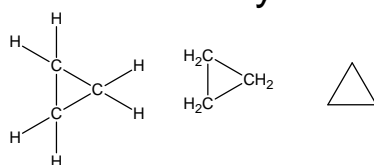


4. If there is more than one methyl group, then each one must have a number locating its carbon atom and a pre-fix of di or tri is given to indicate the total number of methyl groups

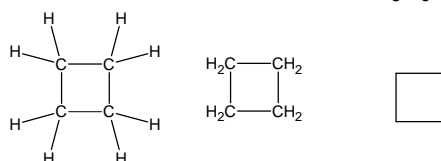


2,2- dimethyl butane

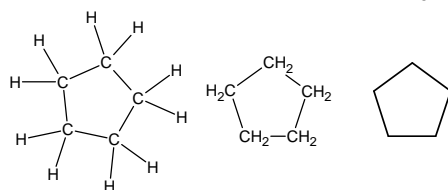
## Hydrocarbon Rings



Cyclopropane. Molecular formula  $C_3H_6$



Cyclobutane. Molecular formula  $C_4H_8$



Cyclopentane. Molecular formula  $C_5H_{10}$

Cycloalkanes  
have the same  
general formula  
as alkenes

e.g.  $C_nH_{2n}$

For **rings** of carbon  
atoms, cyclo- is put  
in front of the name

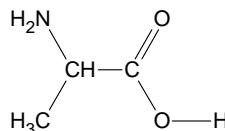
## General naming rules

- The position of the functional group is given by a number, counting from the end that gives the functional group the lowest number (for aldehydes, carboxylic acids & nitriles, the functional group is position 1). *e.g. butanal.*
- Where there are two or more of the same groups, *di-*, *tri-* or *tetra* are used.
- If there is more than one functional group, numbers are separated by commas and the groups are listed in alphabetical order (ignoring *di*, *tri*, etc.). *e.g. 3-bromo-1-chlorobutane, 2,2-dibromo-1-chlorobutane.*
- Numbers are separated by commas and word and numbers by dashes, *e.g. 1-chloro-2,3-dimethylbutane.*
- If a number is not necessary (i.e. the group could only be in one place) then no number should be given. *e.g. propene, methylpropane*

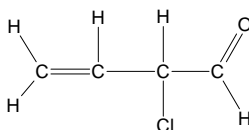
## General naming rules

- Where there are two functional groups, both with suffixes, the preference for the one to have the suffix is carboxylic acid > aldehyde > ketone > alcohol. *e.g. 2-hydroxypropanoic acid, 2-aminopropanoic acid. 2-oxobutanoic acid*

*2-aminopropanoic acid*



- The suffix for alkenes can go in front of other suffixes, *e.g. 2-chlorobut-3-enal.*



## To 'e' or not to 'e'

- When using a suffix, add in the following way :

If the suffix starts with a vowel- remove the -e from the stem alkane name

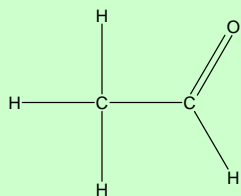
e.g. Propan-1-ol, butan-1-amine, ethanoic acid, ethanoylchloride, butanamide

If the suffix starts with a consonant or there are two or more of a functional group meaning di, or tri needs to be used then **do not remove the the -e** from the stem alkane name

e.g. Propanenitrile, ethane-1,2-diol, propanedioic acid, propane-1,2,3-triol, pentane-2,4-dione

## Naming Carbonyls

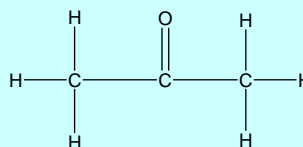
Carbonyls are compounds with a C=O bond, they can be either aldehydes or ketones.



ethanal       $\text{CH}_3\text{CHO}$

If the C=O is on the end of the chain with an H attached it is an aldehyde.

The name will end in -al



$\text{CH}_3\text{COCH}_3$       propanone

If the C=O is in the middle of the chain it is a ketone

The name will end in -one



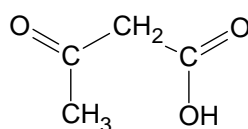
## Naming Carbonyls

Aldehydes never need a number in the name because by definition the functional group is on the end of the chain

Ketones need numbering if there is more than one possibility about where the C=O bond can be.

e.g. propanone and butanone do not need numbers but with 5C's there could be pentan-2-one and pentan-3-one

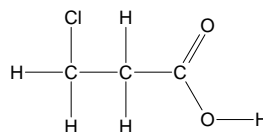
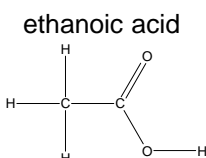
The prefix oxo should be used for compounds that contain a ketone group in addition to a carboxylic acid



3-oxobutanoic acid

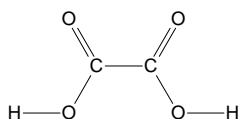
## Naming Carboxylic acid

These have the ending **-oic acid**, but no number is necessary for the acid group as it must always be at the end of the chain. The numbering always starts from the carboxylic acid end.



3-chloropropanoic acid

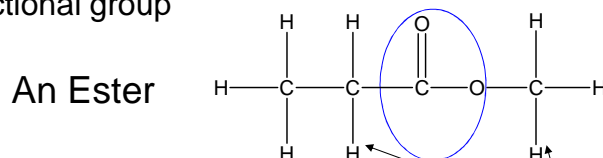
If there are carboxylic acid groups on both ends of the chain then it is called a - dioic acid



Ethanedioic acid  
(CO<sub>2</sub>H)<sub>2</sub>

### Naming Esters

Esters have carbon chains attached to either side of the functional group



Esters have two parts to their names, eg **methyl propanoate**.

The bit ending in **-yl** is next to the single bonded oxygen.

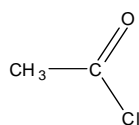
The bit ending in **-anoate** is the chain including the C=O bond

key

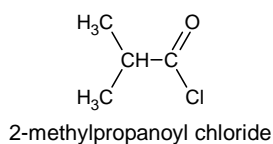
### Naming Acyl Chlorides

Acyl chlorides are carboxylic acid derivatives.

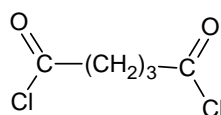
To name take the carboxylic acid name, remove the oic acid bit of the name and replace by **-oyl chloride**



ethanoyl chloride



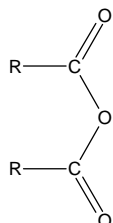
2-methylpropanoyl chloride



Pentanedioyl dichloride

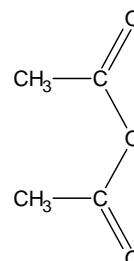
## Naming Acid Anhydrides

Generalised structure of acid anhydride

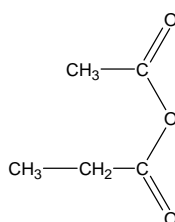


If the R group is the same the name takes the length of R group +1

Ethanoic anhydride



If the groups are different then each one is named e.g. ethanoic propanoic anhydride.



## Naming amines

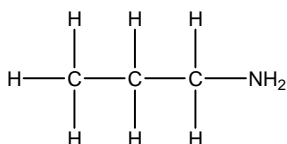
### Amines

These end in **-amine**.

There is, however, rather confusingly two ways of using this suffix.

The exam board tend to use the common version where the name stem ends in **-yl** propylamine.

The IUPAC version of the same chemical is propan-1-amine. (This is used in the same way as naming alcohols)

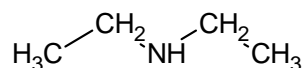


propylamine  
Or propan-1-amine

If the amine is secondary and has two alkyl groups attached to the nitrogen, then each chain is named and the smaller alkyl group is preceded by an -N which plays the same role as a number in positioning a side alkyl chain

$\text{CH}_3\text{CH}_2\text{CH}_2\text{NHCH}_3$   
 N-methylpropylamine  
 (common name)  
 N-methylpropan-1-amine  
 (IUPAC name)

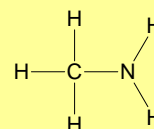
In the common naming version if the chain lengths are the same an -N is not used



Diethylamine (common name- does not use N if chains are same length)  
 N-ethylethanamine (IUPAC name does still use N)

## Naming Amines

primary amine (one C attached to N)

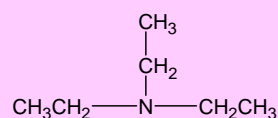


name: methylamine

Secondary amine (two C's attached to N)

$\text{CH}_3\text{CH}_2-\text{NH}-\text{CH}_2\text{CH}_3$   
 name : diethylamine  
 N-ethylethanamine

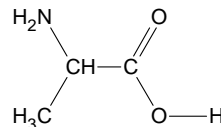
tertiary amine (three C's attached to N)



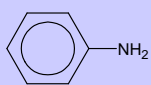
name: triethylamine

## Naming Amines

If there is another functional group as well as the amine group then the prefix amino is used. e.g. 2-aminopropanoic acid.



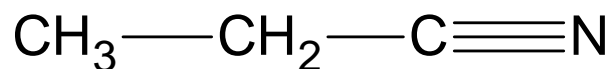
It is also possible to have aromatic amines



phenylamine

## Naming Nitriles

These end in **-ile**, but the C of the CN group counts as one of the chain.



propanenitrile

Note the stem of the name is different to usual (It is propan**en**nitrile and not propan**ni**nitrile)

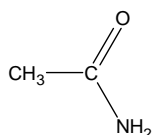
Nitriles never need a number in the name because by the functional group is always on the end of the chain

## Naming Amides

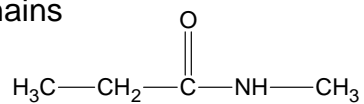
Amides are carboxylic acid derivatives.

Take the carboxylic acid name, remove the oic acid bit of the name and replace by **-amide**

ethanamide



Secondary amides are named differently to show the two carbon chains



N-methyl propanamide