

# IUPAC RULES

## FOR THE NAMING OF ORGANIC COMPOUNDS

- Find the longest chain, which will serve as the basis for the name.
  - If there is a functional group, choose the longest chain that includes the functional group.
- Number the chain from one end to the other.
  - If there is a functional group, start numbering from the end closest to the functional group, meaning the end that assigns the smallest possible number to this group.
  - If there is no functional group, start numbering from the end closest to a substituent, meaning the end that assigns the (first) substituent the smallest number.
- In numerical order, name each substituent according to its position and its type.
  - Group identical substituents together with the Greek prefixes “di-,” “tri-,” “tetra-,” etc.
  - Use commas between numerals, and hyphens to separate numerals from letters.
- After the substituents, name the (most important) functional group and give its numerical position. If several functional groups are present, you must name them all.

Chains:	Hydrocarbon substituents:	Other substituents:	Functional groups:
CH <sub>4</sub> methane	CH <sub>3</sub> - methyl	-F fluoro-	= -ene (double bond)
CH <sub>3</sub> CH <sub>3</sub> ethane	CH <sub>3</sub> CH <sub>2</sub> - ethyl	-Cl chloro-	≡ -yne (triple bond)
CH <sub>3</sub> CH <sub>2</sub> CH <sub>3</sub> propane	CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> - propyl	-Br bromo-	
CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub> butane	(CH <sub>3</sub> ) <sub>2</sub> CH- isopropyl	-I iodo-	-OH -ol (alcohol)
CH <sub>3</sub> (CH <sub>2</sub> ) <sub>3</sub> CH <sub>3</sub> pentane	CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> - butyl	-OH hydroxy-	-CHO -al (aldehyde)
CH <sub>3</sub> (CH <sub>2</sub> ) <sub>4</sub> CH <sub>3</sub> hexane	(CH <sub>3</sub> ) <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> - isobutyl	-NO <sub>2</sub> nitro-	C=O -one (ketone)
CH <sub>3</sub> (CH <sub>2</sub> ) <sub>5</sub> CH <sub>3</sub> heptane	(CH <sub>3</sub> ) <sub>3</sub> C- tert-butyl	-NH <sub>2</sub> amino-	-COOH -oic acid (carboxylic acid)
CH <sub>3</sub> (CH <sub>2</sub> ) <sub>6</sub> CH <sub>3</sub> octane		C=O keto-	
CH <sub>3</sub> (CH <sub>2</sub> ) <sub>7</sub> CH <sub>3</sub> nonane	C <sub>6</sub> H <sub>5</sub> - phenyl		-COO- -oate (ester)
CH <sub>3</sub> (CH <sub>2</sub> ) <sub>8</sub> CH <sub>3</sub> decane			-NH <sub>2</sub> amine
11C is hendecane, etc.	CH <sub>3</sub> O- methoxy,		-CONH <sub>2</sub> amide
20C is eicosane	etc.		-SH thiol
21C is heneicosane, etc.			cyclo- if a ring