#### 2.8

# IUPAC Nomenclature of Unbranched Alkanes

Retained:methane $CH_4$ ethane $CH_3CH_3$ propane $CH_3CH_2CH_3$ butane $CH_3CH_2CH_2CH_3$ 

Note:

*n*-prefix is not part of IUPAC name of any alkane.

For example: *n*-butane is "common name" for CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub>; butane is "IUPAC name."

**Others:** 

Latin or Greek prefix for number of carbons + ane suffix

Number of carbons	Name	Structure
5	pentane	CH <sub>3</sub> (CH <sub>2</sub> ) <sub>3</sub> CH <sub>3</sub>
6	hexane	CH <sub>3</sub> (CH <sub>2</sub> ) <sub>4</sub> CH <sub>3</sub>
7	heptane	CH <sub>3</sub> (CH <sub>2</sub> ) <sub>5</sub> CH <sub>3</sub>
8	octane	CH <sub>3</sub> (CH <sub>2</sub> ) <sub>6</sub> CH <sub>3</sub>
9	nonane	CH <sub>3</sub> (CH <sub>2</sub> ) <sub>7</sub> CH <sub>3</sub>
10	decane	CH <sub>3</sub> (CH <sub>2</sub> ) <sub>8</sub> CH <sub>3</sub>

Number of carbons	Name	Structure
11	undecane	CH <sub>3</sub> (CH <sub>2</sub> ) <sub>9</sub> CH <sub>3</sub>
12	dodecane	$CH_3(CH_2)_{10}CH_3$
13	tridecane	$CH_3(CH_2)_{11}CH_3$
14	tetradecane	$CH_3(CH_2)_{12}CH_3$
15	pentadecane	CH <sub>3</sub> (CH <sub>2</sub> ) <sub>7</sub> CH <sub>3</sub>
16	hexadecane	CH <sub>3</sub> (CH <sub>2</sub> ) <sub>8</sub> CH <sub>3</sub>

Number of carbons Name Structure heptadecane  $CH_3(CH_2)_{15}CH_3$ 17  $CH_{3}(CH_{2})_{16}CH_{3}$ octadecane 18 nonadecane  $CH_3(CH_2)_{17}CH_3$ 19  $CH_{3}(CH_{2})_{18}CH_{3}$ 20 icosane pentacosane CH<sub>3</sub>(CH<sub>2</sub>)<sub>22</sub>CH<sub>3</sub> 25  $CH_{3}(CH_{2})_{28}CH_{3}$ 30 triacontane

# 2.9 Applying the IUPAC Rules: The Names of the $C_6H_{14}$ Isomers





The IUPAC name of the unbranched alkane with a chain of 6 carbons is hexane.

IUPAC Nomenclature of Branched Alkanes (Table 2.7)

Step 1) Find the longest continuous carbon chain and use the IUPAC name of the unbranched alkane as the basis.

Step 2) Add name of substituent as a prefix.

Step 3) Number the chain from the end nearest the substituent, and identify the carbon to which the substituent is attached by number.





according to the number of identical substituents attached to the main chain.

2.10 Alkyl Groups



**Unbranched Alkyl Groups** 

![](_page_14_Picture_1.jpeg)

![](_page_14_Picture_2.jpeg)

If potential point of attachment is at the end of the chain, take the IUPAC name of the corresponding unbranched alkane and replace the *-ane* ending by *-yl*. Unbranched Alkyl Groups

![](_page_15_Picture_1.jpeg)

![](_page_15_Picture_2.jpeg)

CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>

If potential point of attachment is at the end of the chain, take the IUPAC name of the corresponding unbranched alkane and replace the *-ane* ending by *-yl*.

![](_page_15_Figure_4.jpeg)

Unbranched Alkyl Groups

$(\cup \square_2)$	<sub>4</sub> ωΠ <sub>2</sub>	пехуі

 $CH_3(CH_2)_5CH_2$  Heptyl

 $CH_3(CH_2)_{16}CH_2$  Octadecyl

![](_page_17_Figure_1.jpeg)

![](_page_18_Figure_1.jpeg)

IUPAC name: Propyl Common name: *n*-Propyl

### Naming Alkyl Groups (Table 2.8)

Step 1: Identify longest continuous chain starting at point of attachment.

Step 2: Drop -*ane* ending from name of unbranched alkane having same number of carbons as longest continuous chain and replace by -*yl*.

Step 3: Identify substituents on longest continuous chain.

Step 4: Chain is always numbered starting at point of attachment.

![](_page_20_Figure_1.jpeg)

IUPAC name: 1-Methylethyl Common name: Isopropyl

![](_page_21_Figure_1.jpeg)

Classification: Primary alkyl group

Alkyl groups are classified according to the degree of substitution at the carbon that bears the point of attachment. A carbon that is directly attached to one other carbon is a primary carbon.

![](_page_22_Figure_1.jpeg)

Classification: Secondary alkyl group Alkyl groups are classified according to the degree of substitution at the carbon that bears the point of attachment. A carbon that is directly attached to two other carbons is a secondary carbon.

![](_page_23_Figure_1.jpeg)

IUPAC name: Butyl Common name: *n*-Butyl Classification: Primary alkyl group

![](_page_24_Figure_1.jpeg)

IUPAC name: 1-Methylpropyl Common name: sec-Butyl Classification: Secondary alkyl group

![](_page_25_Figure_1.jpeg)

IUPAC name: 2-Methylpropyl Common name: Isobutyl Classification: Primary alkyl group

![](_page_26_Picture_1.jpeg)

IUPAC name: 1,1-Dimethylethyl Common name: *tert*-Butyl Classification: Tertiary alkyl group

# 2.11 IUPAC Names of Highly Branched Alkanes

![](_page_28_Picture_1.jpeg)

![](_page_28_Picture_2.jpeg)

![](_page_29_Picture_1.jpeg)

**4-Ethyloctane** 

![](_page_30_Picture_1.jpeg)

#### 4-Ethyl-3-methyloctane

## List substituents in alphabetical order.

![](_page_31_Picture_1.jpeg)

### 4-Ethyl-3,5-dimethyloctane

# List substituents in alphabetical order. But don't alphabetize di-, tri-, tetra-, etc.

![](_page_32_Figure_0.jpeg)

2,2,6,6,7-Pentamethyloctane?

The chain is numbered in the direction that gives the lower locant to the substituent at the first point of difference in the names. Don't add locants!

#### First Point of Difference Rule

# What is correct name?

#### 2,2,6,6,7-Pentamethyloctane?

The chain is numbered in the direction that gives the lower locant to the substituent at the first point of difference in the names. Don't add locants!

 $C_n H_{2n}$ 

### 2.12

# Cycloalkane Nomenclature

#### Cycloalkanes

Cycloalkanes are alkanes that contain a ring of three or more carbons.

Count the number of carbons in the ring, and add the prefix *cyclo* to the IUPAC name of the unbranched alkane that has that number of carbons.

![](_page_35_Figure_3.jpeg)

**Cycloalkanes** 

Name any alkyl groups on the ring in the usual way.

![](_page_36_Picture_2.jpeg)

Ethylcyclopentane

### Cycloalkanes

Name any alkyl groups on the ring in the usual way.

List substituents in alphabetical order and count in the direction that gives the lowest numerical locant at the first point of difference.

![](_page_37_Figure_3.jpeg)

3-Ethyl-1,1-dimethylcyclohexane