# NOMENCLATURE

(Alkane, Alkenes, Alcohol, Alkyl Halides, Alkynes)

Prepared by Prof. Gade S.T.



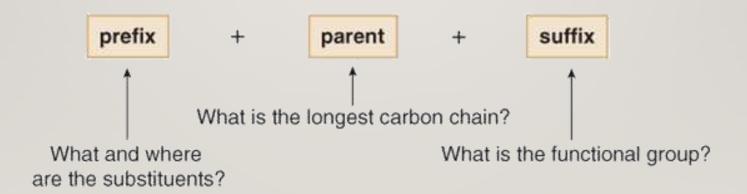
Class: F.Y.B.Pharm

Subject:
Pharmaceutical
Organic Chemistry-I

MULA EDUCATION SOCIETY COLLEGE OF PHARMACY, SONAI

The name of every organic molecule has 3 parts:

- The parent name indicates the number of carbons in the longest continuous chain.
- 2. The **suffix** indicates what **functional group** is present.
- 3. The **prefix** tells us the identity, location, and number of substituents attached to the carbon chain.



# Parent – Longest straight carbon chain

Number of C atoms	Molecular formula	Name (n-alkane)	Number of constitutional isomers
1	CH <sub>4</sub>	methane	_
2	C <sub>2</sub> H <sub>6</sub>	ethane	_
3	C <sub>3</sub> H <sub>8</sub>	propane	_
4	C <sub>4</sub> H <sub>10</sub>	butane	2
5	C <sub>5</sub> H <sub>12</sub>	pentane	3
6	C <sub>6</sub> H <sub>14</sub>	hexane	5
7	C <sub>7</sub> H <sub>16</sub>	heptane	9
8	C <sub>8</sub> H <sub>18</sub>	octane	18
9	C <sub>9</sub> H <sub>20</sub>	nonane	35
10	C <sub>10</sub> H <sub>22</sub>	decane	75
20	C <sub>20</sub> H <sub>42</sub>	eicosane	366,319

**Suffix** — Our first functional group is alkane, so the suffix is —ane For later functional groups we will drop the —ane root suffix for others

Alkane chain	# Carbons	Name
CH <sub>4</sub>	1	methane
CH <sub>3</sub> CH <sub>3</sub>	2	ethane
CH <sub>3</sub> CH <sub>2</sub> CH <sub>3</sub>	3	propane
CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>	4	butane
CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>	5	pentane
CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>	6	hexane
CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>	7	heptane
CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>	8	octane
CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>	9	nonane
CH <sub>3</sub> CH <sub>2</sub> CH <sub>3</sub>	10	decane
CH <sub>3</sub> CH <sub>2</sub>	11	undecane
CH <sub>3</sub> CH <sub>2</sub>	12	dodecane

Prefix — Our substituents will be branches in the alkane structure

A branch is another alkane minus one hydrogen – an alkyl group

**Example** – if CH<sub>3</sub>- is a branch on a longer chain:

CH<sub>3</sub>- is CH<sub>4</sub> minus 1 hydrogen

Since it is a side chain we will replace the –ane suffix with –yl

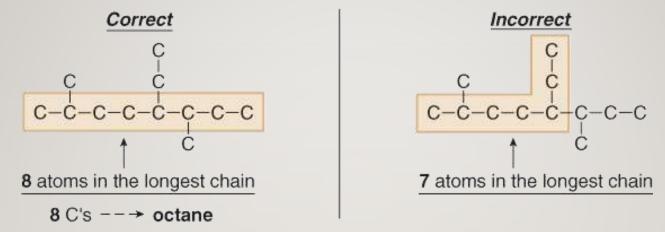
CH<sub>3</sub>- is a methyl group

We can also abbreviate this group as Me-

# **Prefixes** -

Alkyl group	Structure	IUPAC name	Abbreviation
CH <sub>3</sub> -	CH <sub>3</sub> -	methyl	Me-
CH <sub>3</sub> CH <sub>2</sub> -	Ser s	ethyl	Et-
CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> -	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	<i>n</i> -propyl	<i>n</i> -Pr
CH₃CHCH₃ 	- Sept	isopropyl or <i>i</i> -propyl	<i>i</i> -Pr
CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> -	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	<i>n</i> -butyl	<i>n</i> -Bu
CH <sub>3</sub> CH <sub>2</sub> CHCH <sub>3</sub>	- Ser	<i>sec</i> -butyl	<i>s</i> -Bu
(CH <sub>3</sub> ) <sub>2</sub> CHCH <sub>2</sub> -	Sec. 2	isobutyl or <i>i</i> -butyl	<i>i</i> -Bu
(CH <sub>3</sub> ) <sub>3</sub> C-	× <sub>s</sub> x³	tert-butyl or t-butyl	<i>t</i> -Bu
C <sub>6</sub> H <sub>5</sub> -		phenyl	Ph

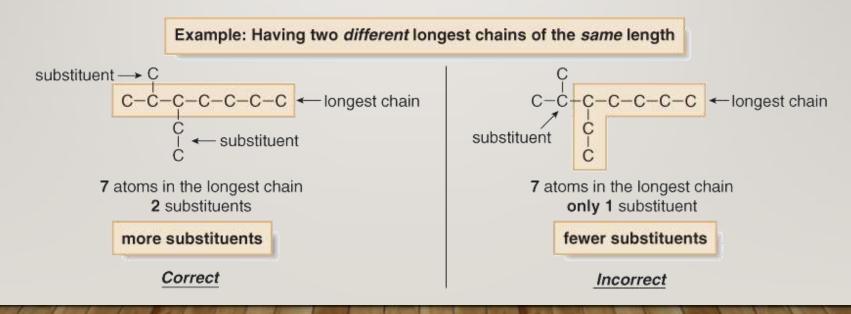
# 1. Find the parent carbon chain and add the suffix.



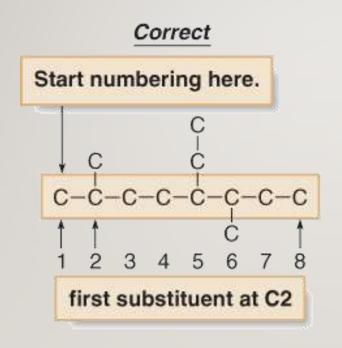
Note that it does not matter if the chain is straight or it bends.

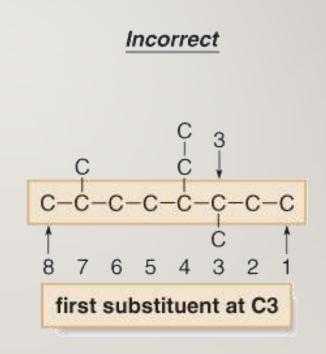
Also note that if there are two chains of equal length, pick the chain with *more* substituents.

In the following example, two different chains in the same alkane have seven C atoms. We circle the longest continuous chain as shown in the diagram on the left, since this results in the greater number of substituents.



2. Number the atoms in the carbon chain to give the **first substituent** the **lowest number**.

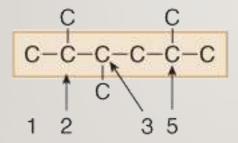




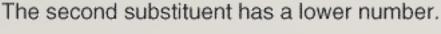
If the first substituent is the same distance from both ends, number the chain to give the second substituent the lower number.

## Example: Giving a lower number to the second substituent

### Numbering from left to right

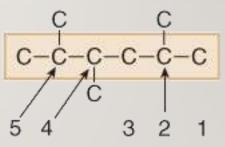


CH<sub>3</sub> groups at C2, C3, and C5.



Correct

### Numbering from right to left



CH<sub>3</sub> groups at C2, C4, and C5.

higher number

When numbering a carbon chain results in the same numbers from either end of the chain, assign the lower number alphabetically to the first substituent.

Example: Two different groups equidistant from the ends

### Numbering from left to right

- ethyl at C3
- methyl at C5

Earlier letter → lower number

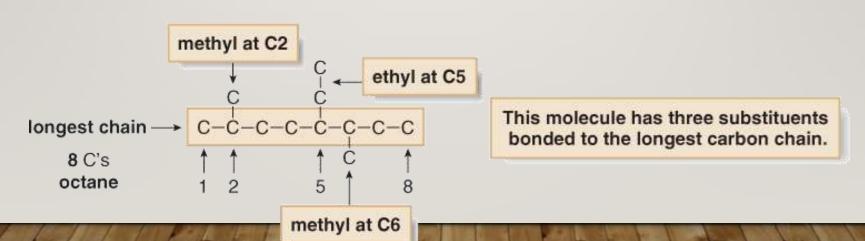
Correct

### Numbering from right to left

- methyl at C3
- ethyl at C5

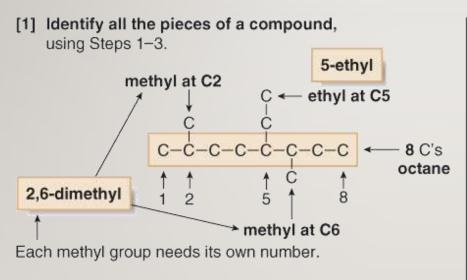
Incorrect

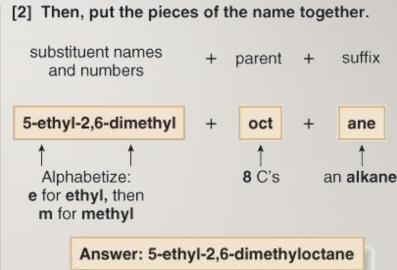
- 3. Name and number the substituents
  - Name the substituents as alkyl groups.
  - Every carbon belongs to either the longest chain or a substituent, not both.
  - Each substituent needs its own number
  - If two or more identical substituents are bonded to the longest chain, use prefixes to indicate how many: di- for two groups, trifor three groups, tetra- for four groups, and so forth.



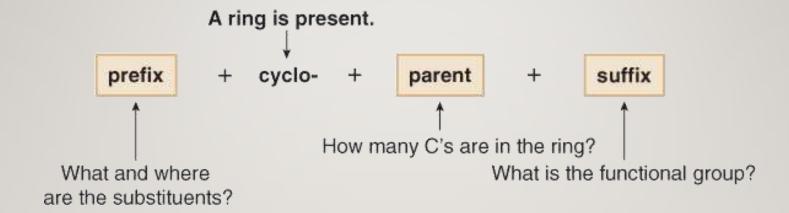
- 4. Combine substituent names and numbers + parent and suffix.
  - Precede the name of the parent by the names of the substituents.
  - Alphabetize the names of the substituents, ignoring all prefixes except iso, as in isopropyl and isobutyl.
  - Precede the name of each substituent by the number that indicates its location.
  - Separate numbers by commas and separate numbers from letters by hyphens.
  - The name of an alkane is a single word, with no spaces after hyphens and commas.

4. Combine substituent names and numbers + parent and suffix.





**Cycloalkanes** are named by using similar rules, but the prefix cyclo- immediately precedes the name of the parent.



1. Find the parent cycloalkane.

2. Name and number the substituents. No number is needed to indicate the location of a single substituent.

For rings with more than one substituent, begin numbering at one substituent and proceed around the ring to give the second substituent the lowest number.

### numbering clockwise

CH<sub>3</sub> groups at C1 and **C3**The 2<sup>nd</sup> substituent has a lower number.

Correct: 1,3-dimethylcyclohexane

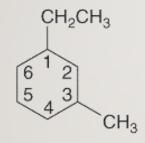
### numbering counterclockwise

CH3 groups at C1 and C5

Incorrect: 1,5-dimethylcyclohexane

With two different substituents, number the ring to assign the lower number to the substituents **alphabetically**.

Begin numbering at the ethyl group.

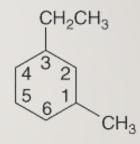


- ethyl group at C1
- methyl group at C3

earlier letter → lower number

Correct: 1-ethyl-3-methylcyclohexane

Begin numbering at the methyl group.

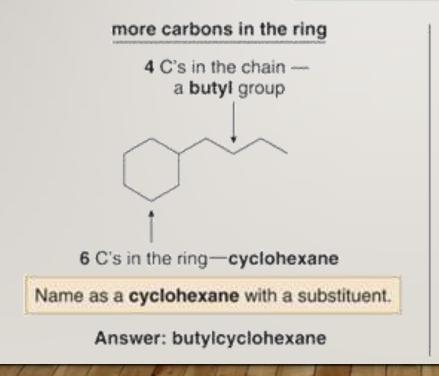


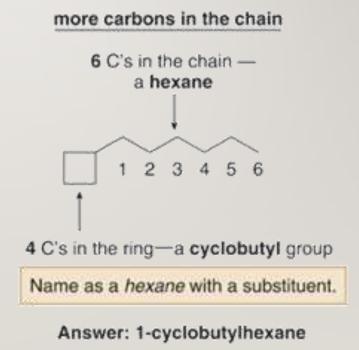
- methyl group at C1
- ethyl group at C3

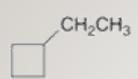
Incorrect: 3-ethyl-1-methylcyclohexane

Note the special case of an alkane composed of both a ring and a long chain. If the number of carbons in the ring is greater than or equal to the number of carbons in the longest chain, the compound is named as a cycloalkane.

Contrast two different examples





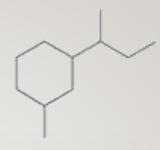


### ethylcyclobutane

No number is needed with only one substituent.

### 1,2-dimethylcyclohexane

Number to give the 2<sup>nd</sup> CH<sub>3</sub> group the lower number: 1,2- not 1,6-.



### 1-sec-butyl-3-methylcyclohexane

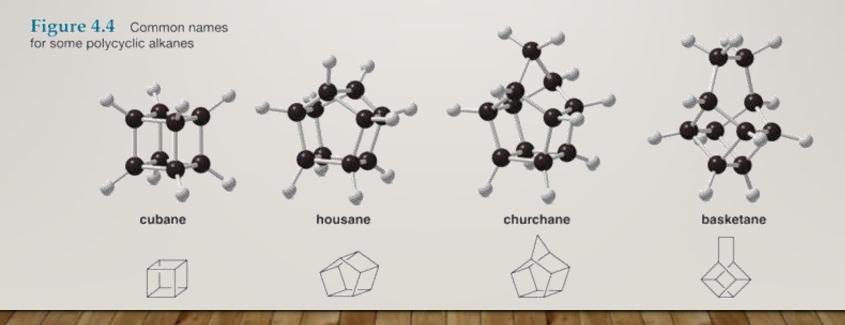
Assign the lower number to the 1st substituent alphabetically: the **b** of **b**utyl before the **m** of **m**ethyl.

### 1,2,4-triethylcyclopentane

Number to give the 2<sup>nd</sup> CH<sub>3</sub>CH<sub>2</sub> group the lower number: 1,2,4- not 1,3,4- or 1,3,5-.

### Nomenclature—Common Names

Some organic compounds are identified using common names that do not follow the IUPAC system of nomenclature. Many of these names were given long ago before the IUPAC system was adopted, and are still widely used. Additionally, some names are descriptive of shape and structure, like those below:



# **Other Functional Groups - Nomenclature**

The IUPAC rules for all other functional groups will differ only by the following:

- 1. Suffix will change to reflect functional group
- 2. Some functional groups have priority over others
- 3. For now: alcohol > alkyne > alkene > alkane=alkyl halide
- 4. The longest chain must contain the suffix functional group even if not the longest chain overall
- 5. Numbering gives this functional group the lowest number even if there are other groups that would be lower
- 6. If an alkene has <u>stereochemistry</u>, it must be specified in the prefix

# **Alkyl Halides - Nomenclature**

### HOW TO Name an Alkyl Halide Using the IUPAC System

Example Give the IUPAC name of the following alkyl halide:

Step [1] Find the parent carbon chain containing the halogen.

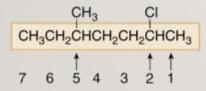
7 C's in the longest chain

7 C's ---→ heptane

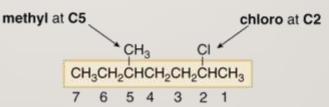
 Name the parent chain as an alkane, with the halogen as a substituent bonded to the longest chain.

### Step [2] Apply all other rules of nomenclature.

a. Number the chain.



 Begin at the end nearest the first substituent, either alkyl or halogen. b. Name and number the substituents.



c. Alphabetize: c for chloro, then m for methyl.

ANSWER: 2-chloro-5-methylheptane

# **Alcohols - Nomenclature**

### HOW TO Name an Alcohol Using the IUPAC System

**Example** Give the IUPAC name of the following alcohol:

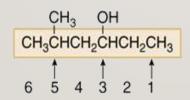
Step [1] Find the longest carbon chain containing the carbon bonded to the OH group.

6 C's -→ hexane -→ hexanol

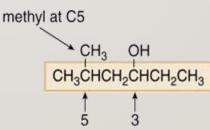
 Change the -e ending of the parent alkane to the suffix -ol.

Step [2] Number the carbon chain to give the OH group the lower number, and apply all other rules of nomenclature.

a. Number the chain.



 Number the chain to put the OH group at C3, not C4. b. Name and number the substituents.



Answer: 5-methyl-3-hexanol

3-hexanol

# **Alcohols - Nomenclature**

- When an OH group is bonded to a ring, the ring is numbered beginning with the OH group.
- Because the functional group is at C1, the 1 is usually omitted from the name.
- The ring is then numbered in a clockwise or counterclockwise fashion to give the next substituent the lowest number.

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

3-methylcyclohexanol

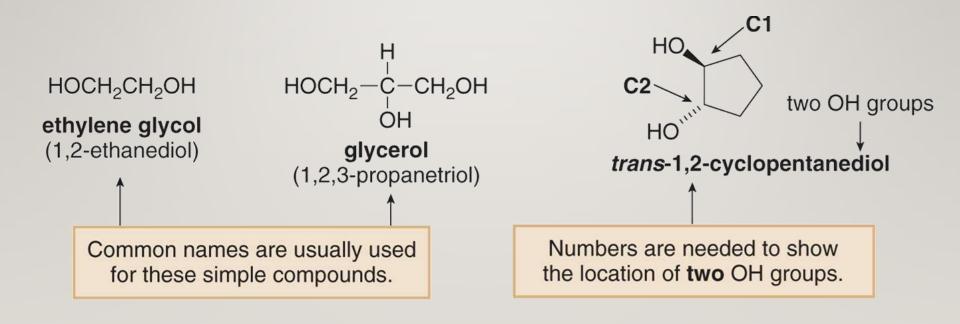
2,5,5-trimethylcyclohexanol

The OH group is at C1; the second substituent (CH<sub>3</sub>) gets the lower number.

The OH group is at C1; the second substituent (CH<sub>3</sub>) gets the lower number.

# **Alcohols - Nomenclature**

- Compounds with two hydroxyl groups are called diols or glycols.
- Compounds with three hydroxyl groups are called triols.



HOW TO Name an Alkene

Example Give the IUPAC name of the following alkene:

$$CH_3$$
  $CH_3$   $CH_2CHCH_2CHCH_3$ 

Step [1] Find the longest chain that contains both carbon atoms of the double bond.

6 C's in the longest chain

hexane ---→ hexene

• Change the -ane ending of the parent alkane to -ene.

**Step [2]** Number the carbon chain to give the double bond the lower number, and apply all other rules of nomenclature.

a. **Number** the chain, and name using the *first* number assigned to the C=C.

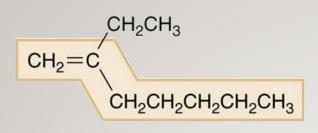
 Number the chain to put the C=C at C2, not C4.

2-hexene

b. Name and number the substituents.

three methyl groups at C2, C3, and C5

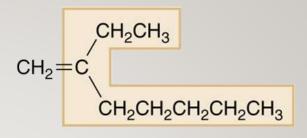
Answer: 2,3,5-trimethyl-2-hexene



7 C's ---→ heptene

Both C's of the C=C are contained in this long chain.

Correct: 2-ethyl-1-heptene

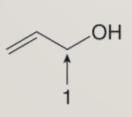


8 C's

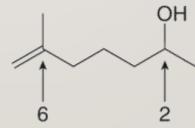
Both C's of the C=C are NOT contained in this long chain.

Incorrect

 Compounds that contain both a double bond and a hydroxy group are named as alkenols and the chain (or ring) is numbered to give the OH group the lower number.

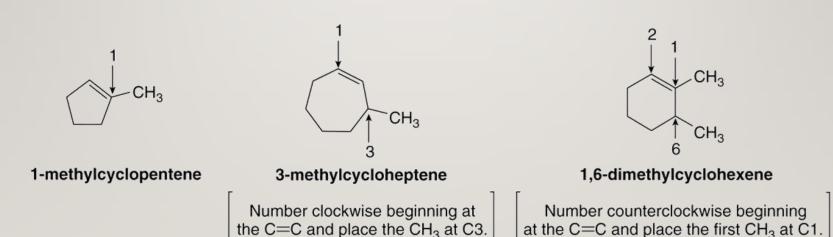


2-propen-1-ol



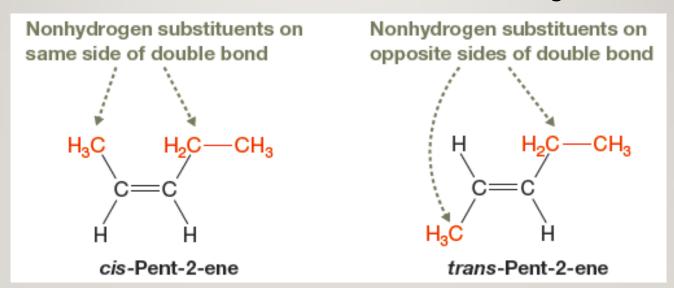
6-methyl-6-hepten-2-ol

- Compounds with two double bonds are named as dienes by changing the "-ane" ending of the parent alkane to the suffix "-adiene".
- Compounds with three double bonds are named as trienes, and so forth.
- In naming cycloalkenes, the double bond is located between C1 and C2, and the "1" is usually omitted in the name.
- The ring is numbered clockwise or counterclockwise to give the first substituent the lower number.



# STEREOCHEMICAL CONFIGURATIONS OF ALKENES: Z/E DESIGNATIONS

There are limitations to cis and trans alkene designations.

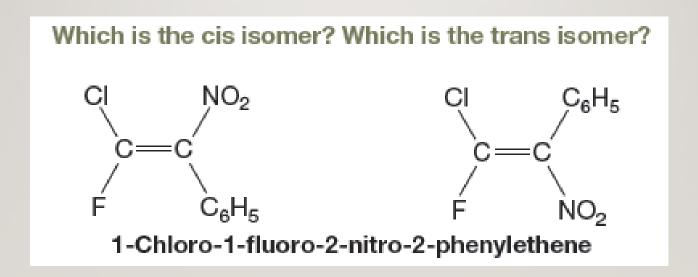


- A cis alkene describes two nonhydrogen substituents that are on the same side of the double bond.
- A trans alkene describes two nonhydrogen substituents that are on opposite sides of the double bond.

# STEREOCHEMICAL CONFIGURATIONS OF ALKENES: Z/E DESIGNATIONS

CONTINUED...

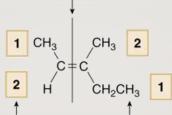
 When there is more than one nonhydrogen substituent bonded to one or both alkene carbons, you cannot use cis or trans designations.



### HOW TO Assign the Prefixes E and Z to an Alkene

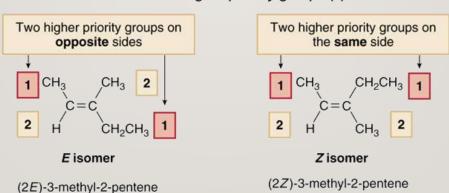
- **Step** [1] Assign priorities to the two substituents on each end of the C=C by using the priority rules for *R*,*S* nomenclature (Section 5.6).
  - Divide the double bond in half, and assign the numbers 1 and 2 to indicate the relative priority of the two groups on each end—the higher priority group is labeled 1, and the lower priority group is labeled 2.

Divide the double bond in half.



Assign priorities to each side of the C=C separately.

### **Step [2]** Assign E or Z based on the location of the two higher priority groups (1).



- The E isomer has the two higher priority groups on the opposite sides.
- The **Z** isomer has the two higher priority groups on the same side.

- Alkynes are named in the same general way that alkenes are named.
- In the IUPAC system, change the –ane ending of the parent alkane name to the suffix –yne.
- Choose the longest continuous chain that contains both atoms of the triple bond and number the chain to give the triple bond the lower number.

- Compounds with two triple bonds are named as diynes, those with three are named as triynes and so forth.
- Compounds with both a double and triple bond are named as enynes.
- The chain is numbered to give the first site of unsaturation (either C=C or C≡C) the lower number.

# THANK YOU