

Recall:**Nomenclature of Alkene**

- Note that there is no free rotation around $C=C$.

Rules:

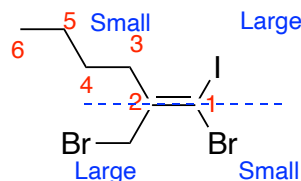
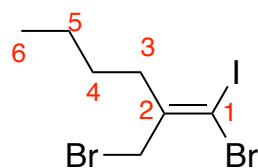
1. Find longest chain
2. Number from end to contain both ends of $C=C$ and give lowest number to 1st C of $C=C$
3. Change “ane” to “ene” precede with number to indicate first double bond position

E/Z Nomenclature

E - Entgegen - Opposite

Z - Zusammen - Together

Naming based on atomic number, similar process to identifying S/R stereochemistry

Example:

Substituents: 1-iodo; 1-bromo; 2-(1-bromomethyl)
Parent: Hex-1-ene or 1-Hexene

Name: *E*-1-bromo-2-(1-bromomethyl)-1-iodo-1-hexene

Nomenclature of Cycloalkenes

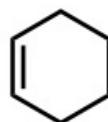
Cyclopropene



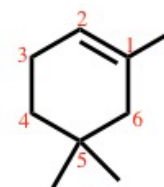
Cyclobutene



Cyclopentene

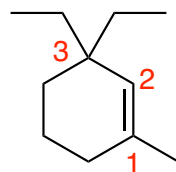


Cyclohexene



1,5,5-trimethyl-1-cyclohexene

Rule: Number the cycloalkene such that the double bond is between C1 and C2 and that the first substituent has the lowest number possible.

Example:

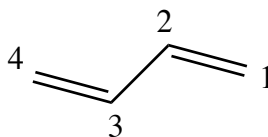
3,3-diethyl-1-methyl-1-cyclohexene or
3,3-diethyl-1-methylcyclohex-1-ene

Nomenclature of alkenes with multiple carbon-carbon double bonds (poly-enes):Multiple $\text{C}=\text{C}$

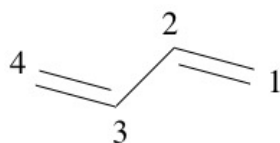
2 Diene

3 Triene

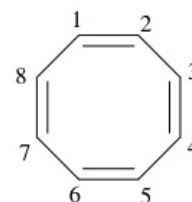
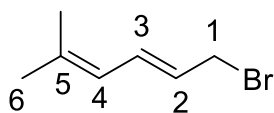
4 Tetraene

↓
...etcButa-1,3-Diene
1,3-ButadieneDrop -ne and add "diene",
"triene", etc.

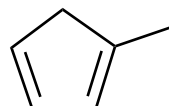
- 1) Find the longest chain containing the maximum number of double bonds.
- 2) Start numbering such that the first doubly bonded position would have the lowest number possible
- 3) Write out the full name. Number the substituents according to their position in the chain and list them alphabetically.

Buta-1,3-diene
1,3-Butadiene

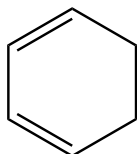
1,3-Cyclobutadiene

1,3,5,7-Cyclooctatetraene (All Z)
COT

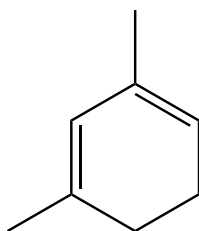
3E-1-bromo-5-methyl-2,4-hexadiene



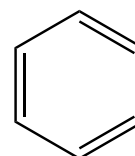
1-methyl-1,3-cyclopentadiene

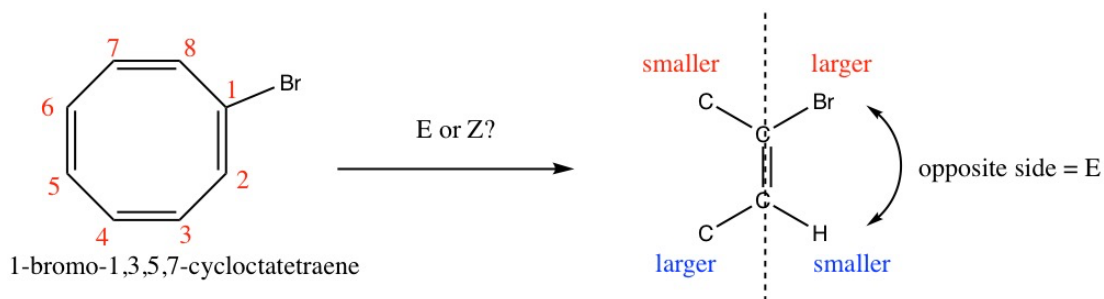
Other examples:

1,3-cyclohexadiene

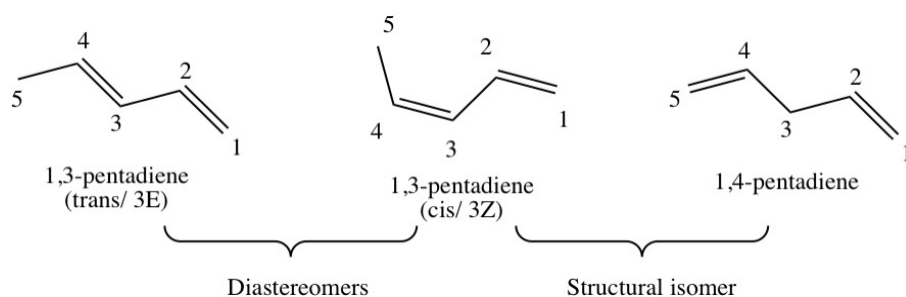


1,3-dimethyl-1,3-cyclohexadiene

benzene
(NOT a cyclohexatriene)
(aromatic)Abbrev:
PhH or C_6H_6 or ϕH

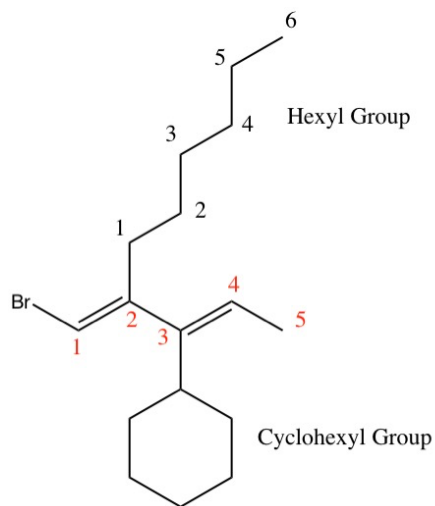


It is therefore (E)-1-bromo-1,3,5,7-cyclooctatetraene



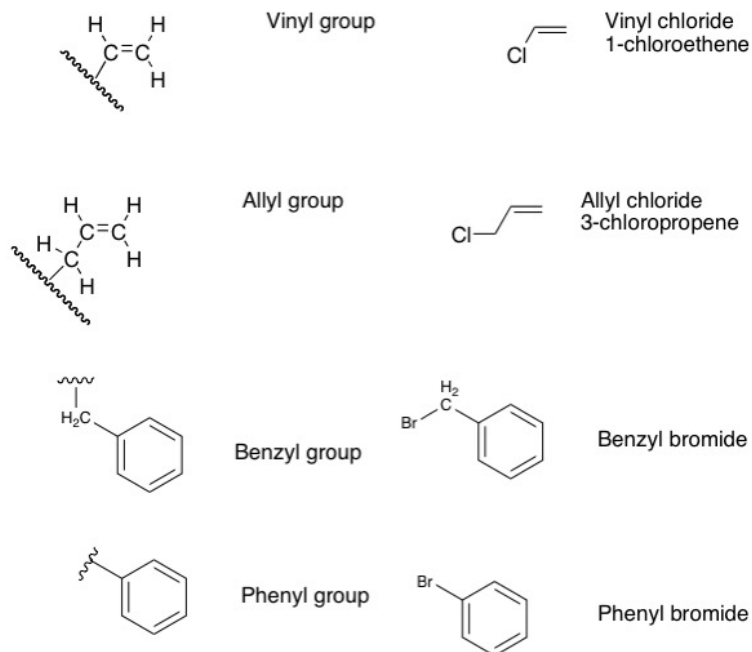
1,3-pentadiene (trans) = (E)-1,3-pentadiene

1,3-pentadiene (cis) = (Z)-1,3-pentadiene



1E,3E-1-Bromo-3-cyclohexyl-2-hexyl-1,3-pentadiene

Note: Carbons attached to double and triple bonds are depicted as additional carbon-carbon bonds in the representations above.

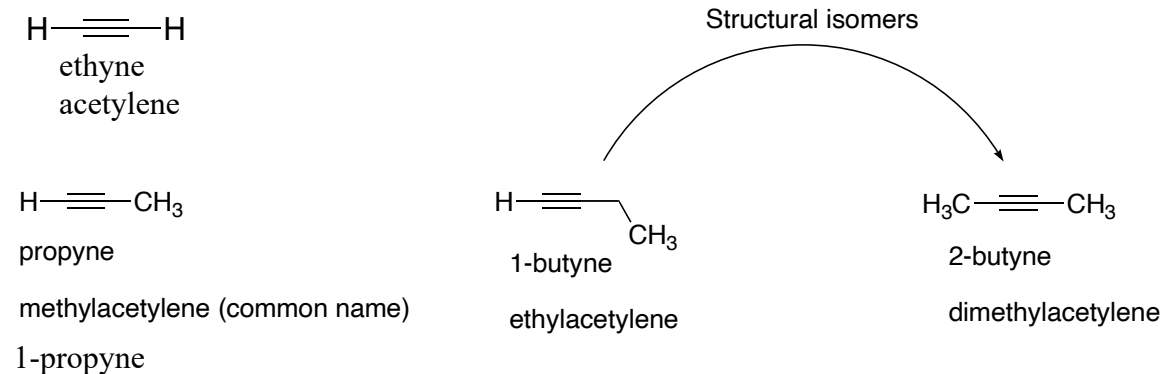
Special Nomenclature of Common Groups:

Note: phenyl bromide is commonly called bromobenzene

Nomenclature of Alkynes (also known as acetylenes)

Rules:

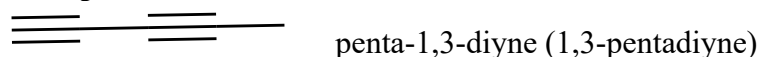
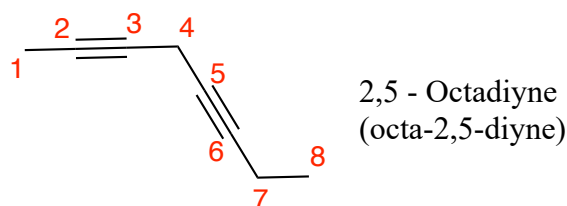
- Find longest chain with max number of multiple bonds
- Number from end to give 1st multiply bonded position the lowest number
- Drop “ane” and add “yne”
- For multiple triple bonds, drop “ne” and add “diyne”, “triyne”, etc.
- Halides and alkyl substituents take lower priority than double or triple bond



Multiple alkynes end with:

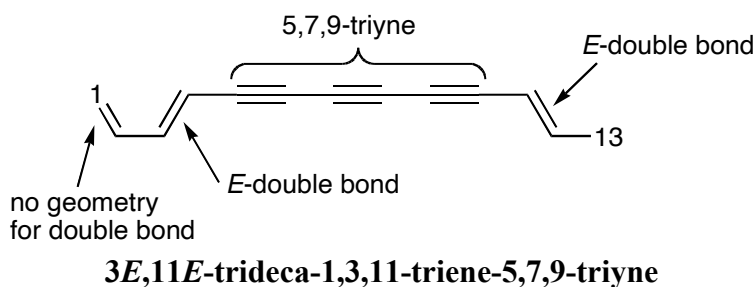
- 2 $\text{C} \equiv \text{C}$ diyne
 3 $\text{C} \equiv \text{C}$ triyne
 4 $\text{C} \equiv \text{C}$ tetrayne

Mixed double and triple bond containing compounds are “eneynes.”

Example 1:**Example 2****Example 3:****Example 4:**

The below example is from canola – defense substance (anti-nematode)

- Parent alkane of 13 carbons is tridecane – hence trideca
- Start numbering the chain such that the **first multiply bonded position** gets the lowest number possible.



Note: alkene stereochemistry can go right to the numbers indicating positions of double bonds:
trideca-1,3*E*,11*E*-triene-5,7,9-triyne

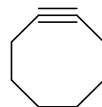
Cyclic Structures with Triple Bonds



extremely unstable
Cyclohexyne



$T_{1/2}$ approx 1 h
very reactive
Cycloheptyne



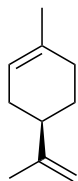
Cyclooctyne

- Cycloalkynes are very reactive as the triple bonds wants to assume a 180° bond angle.
- Cycloheptyne is more stable as the angle is larger ($T_{1/2}=1$ h)
- Cyclooctyne is known to be relatively stable compound used in azide-alkyne [3+2] cycloadditions (used in biorthogonal coupling – Nobel Prize 2022 Carolyn Bertozzi & Barry Sharpless)

Terpenes or Isoprenoids or Terpenoids

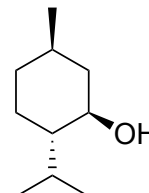


isoprene
2-methyl-1,3-butadiene
 C_5 unit

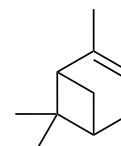


limonene
citrus
 C_{10} (two C_5 units)

- 1 C_5 Hemiterpene
- 2 C_5 Monoterpene
- 3 C_5 Sesquiterpene
- 4 C_5 Diterpene
- 5 C_5 Sesterterpene
- 6 C_5 Triterpene



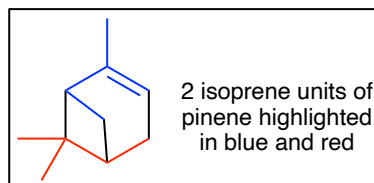
menthol
monoterpene



pinene
monoterpene

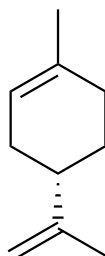
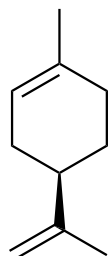
- Steroids are made from triterpenes

- Pinene is made from two isoprene units

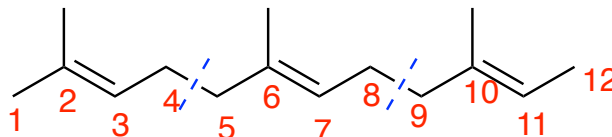


- Leopold Ruzicka: did significant work with terpenes

Examples:



Enantiomers of Limonene



2,6,10-trimethyldodeca-2,6,10-triene

2,6,10-trimethyl-2,6,10-dodecatriene