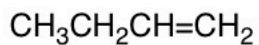


REVIEW:**Alkene Nomenclature:**

butylene

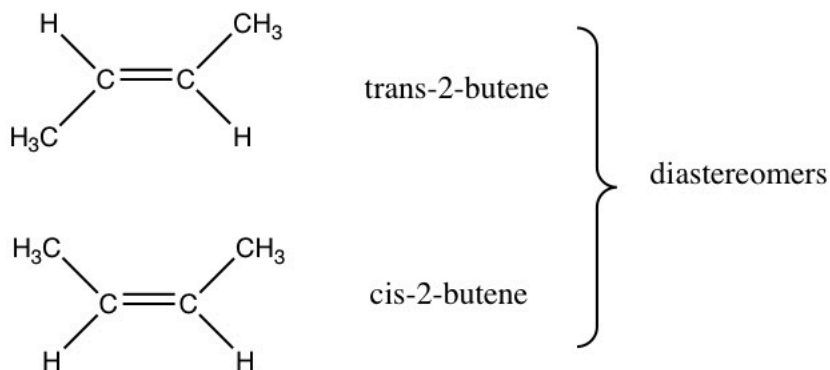
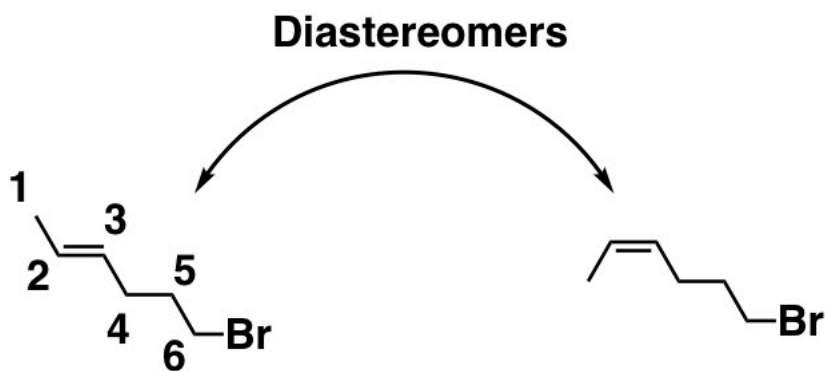
OR

1-butene OR



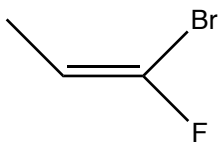
but-1-ene

Below are two structural isomers of 1-butene

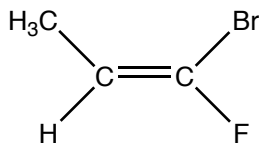
**Example 1: 6-Bromo-2-hexene****trans-6-Bromo-2-hexene****cis-6-Bromo-2-hexene**

In the cis isomer, the two higher priority groups on either side of the carbon-carbon double bond are pointing in the same direction.

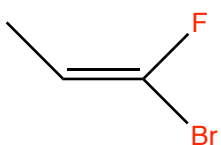
Example 2: 1-Bromo-1-fluoro-1-propene



means



1-bromo-1-fluoropropene



1-bromo-1-fluoropropene

Question: Are the compounds above the same?

Answer: No, they are diastereomers and we can differentiate them by using the E and Z nomenclature

E, Z - Nomenclature

E - Entgegen - Opposite

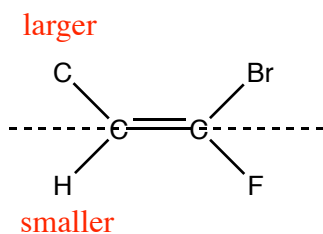
Z - Zusammen - Together

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

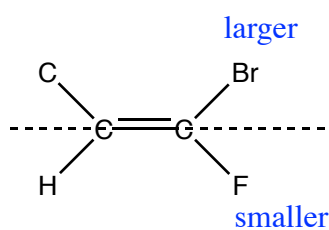
Example 1: 1-bromo-1-fluoro-1-propene

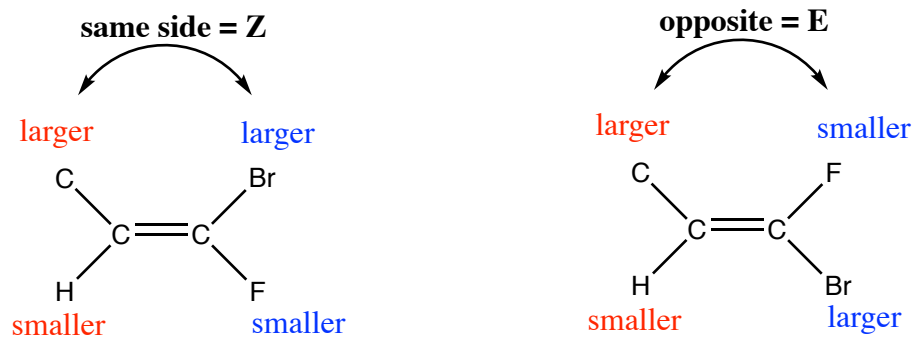
- compare the atomic no. of the adjacent atoms.

Compare the **left** side of the C=C bond

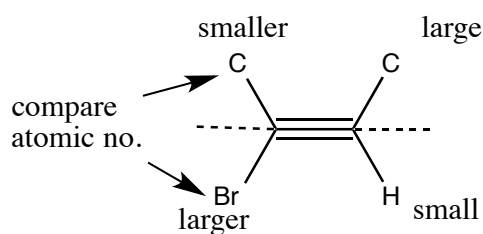


Compare the **right** side of the C=C bond

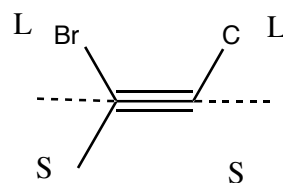




Example 2: 2-bromo-2-butene

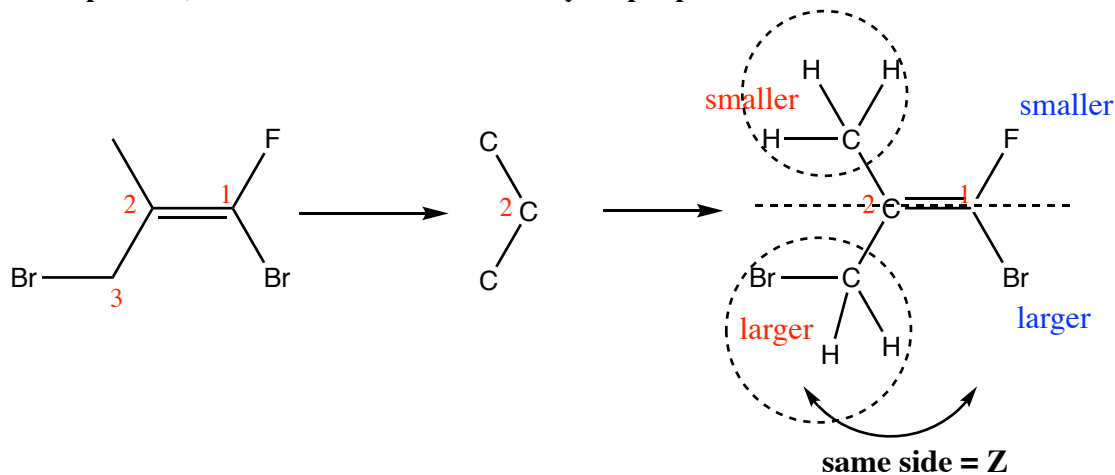


Large groups are on opposite sides on the C=C --> E
E-2-bromo-2-butene



Z-2-bromo-2-butene

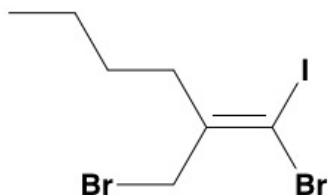
Example 3: 1,3-dibromo-1-fluoro-2-methyl-1-propene



Therefore the name is: (Z)-1,3-dibromo-1-fluoro-2-methyl-1-propene

If you cannot decide on basis of atomic number of atoms directly attached to double bond, go to the next set of atoms until a higher atomic number is found

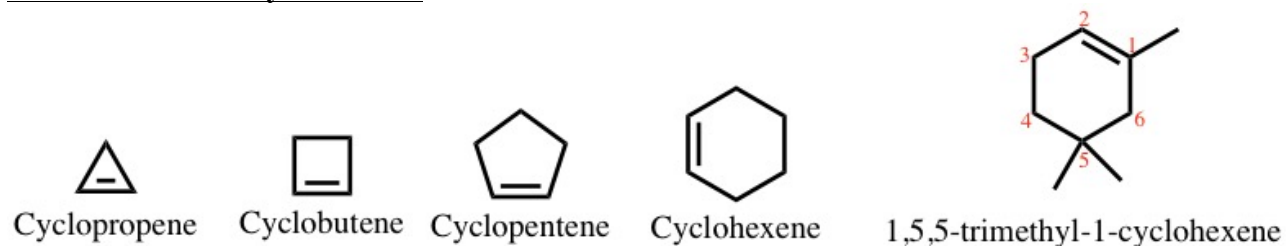
Example 3:



E-1-Bromo- 2-bromomethyl-1-iodohex-1-ene

Iodine is on the opposite side to the bromomethyl (highest priority groups on either side of the alkene) and so the stereochemistry is deemed E.

Nomenclature of Cycloalkenes



Nomenclature of alkenes with multiple carbon-carbon double bonds (poly-enes):

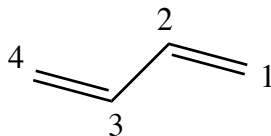
Multiple $C=C$

2 Diene

3 Triene

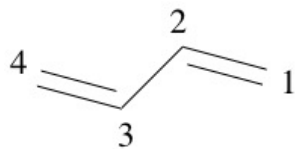
4 Tetraene

...etc



Buta-1,3-Diene
1,3-Butadiene

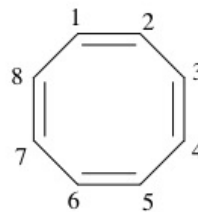
Drop -ne and add “diene”,
“triene”, etc.



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

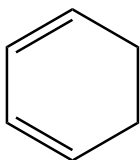


1,3-Cyclobutadiene

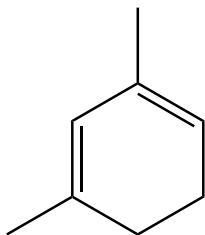


1,3,5,7-Cyclooctatetraene
COT

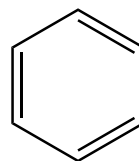
Other examples:



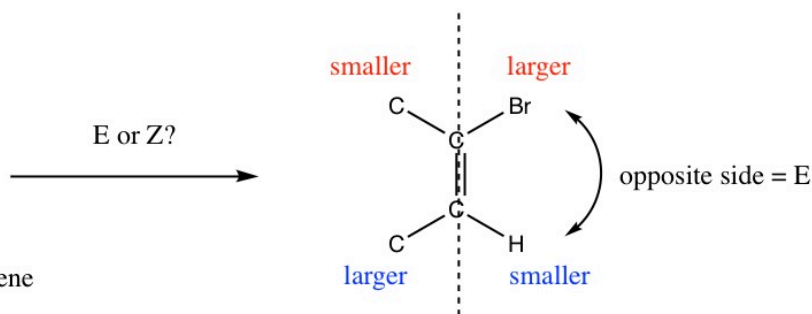
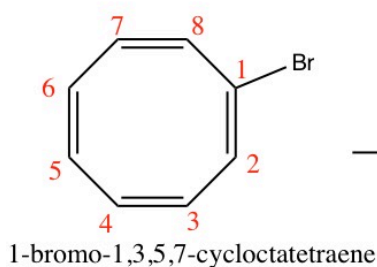
1,3-cyclohexadiene



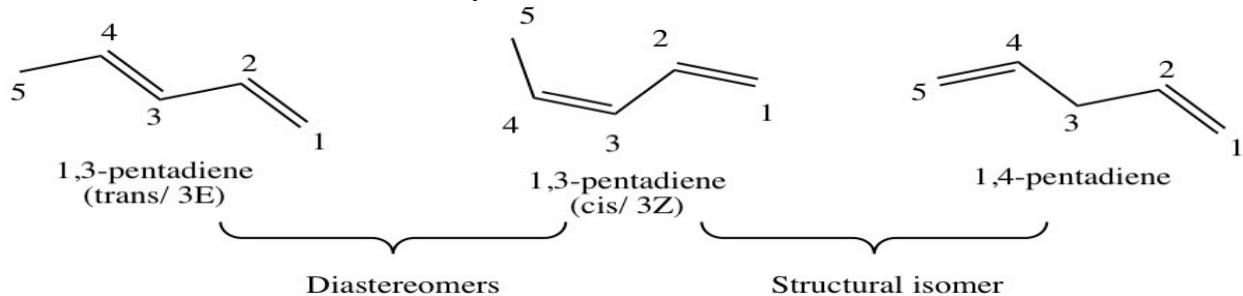
1,3-dimethyl-1,3-cyclohexadiene



benzene
(NOT a cyclohexatriene)
(aromatic)

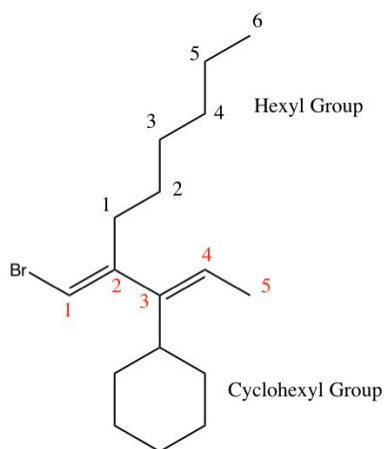


It is therefore (E)-1-bromo-1,3,5,7-cyclooctatetraene



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

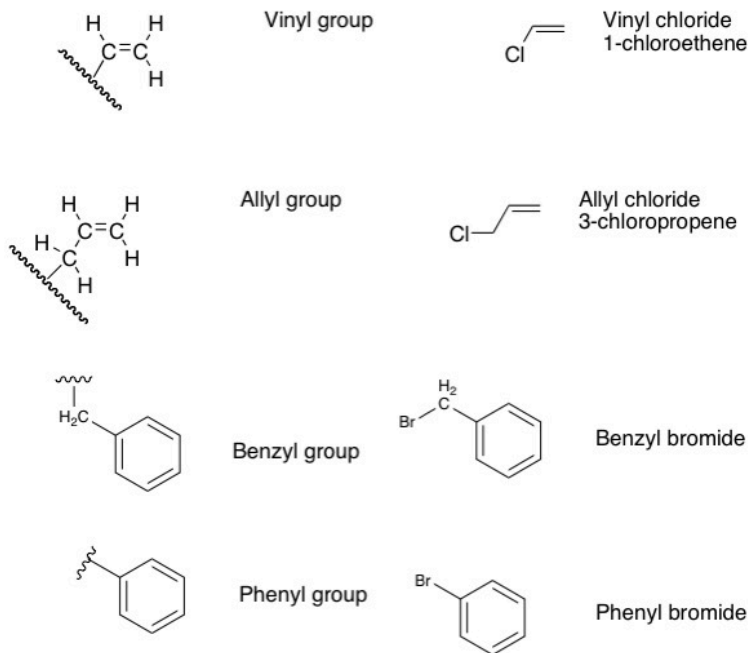
1,3-pentadiene (cis) = (Z)-1,3-pentadiene



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:

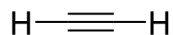


phenyl bromide is commonly called bromobenzene

Nomenclature of Alkynes

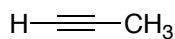
Rules:

- 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.



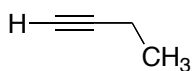
ethyne
acetylene

Structural isomers



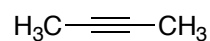
propyne

methylacetylene (common name)



1-butyne

ethylacetylene



2-butyne

dimethylacetylene

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:

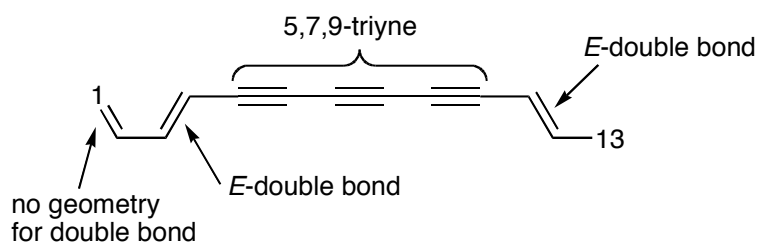


2,4 - Octadiyne

Example 2:

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

Parent alkane of 13 carbons is tridecane – hence trideca



3E,11E-trideca-1,3,11-triene-5,7,9-triyne