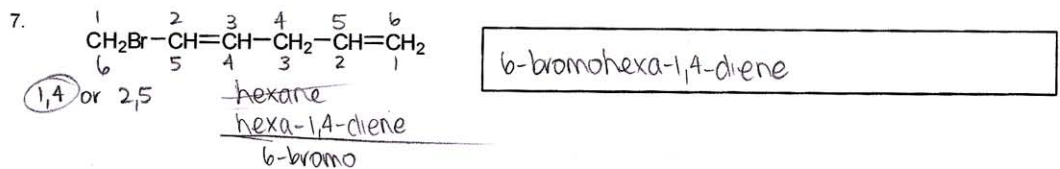
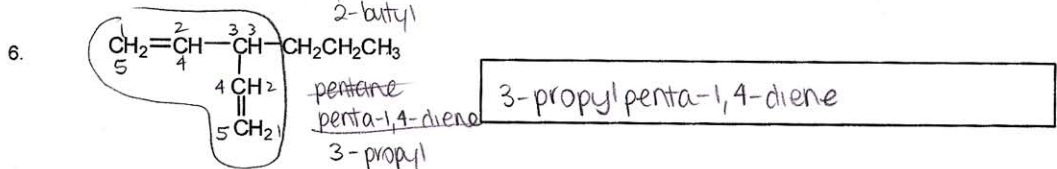
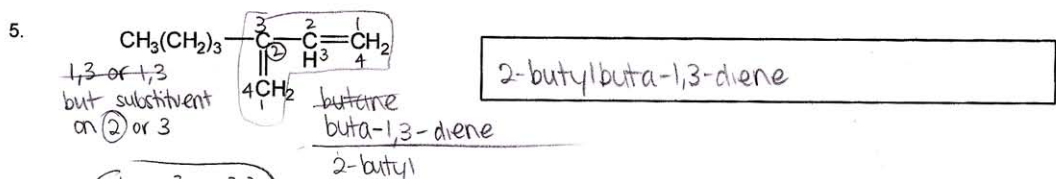
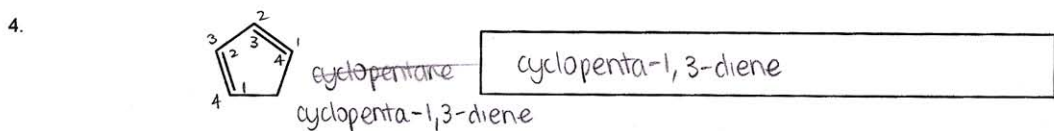
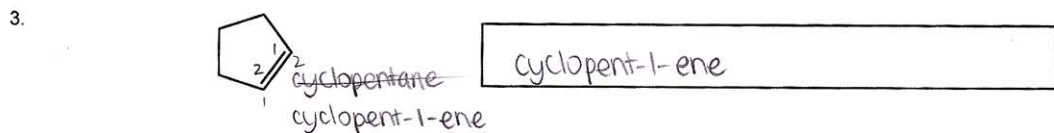
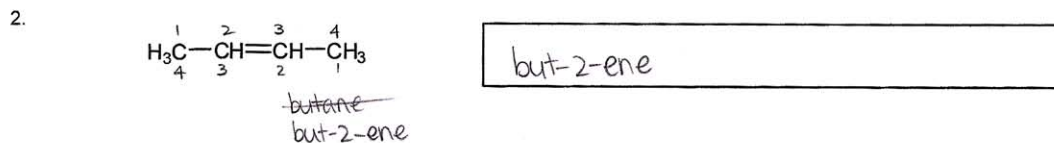
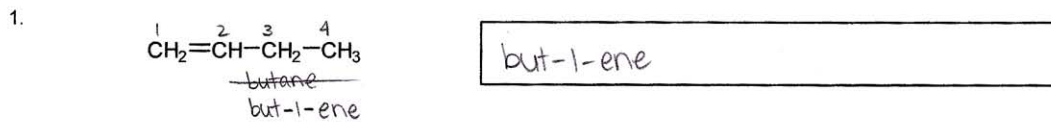
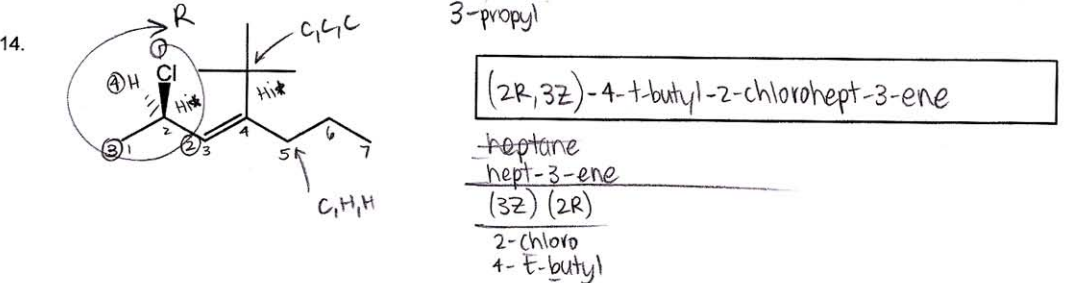
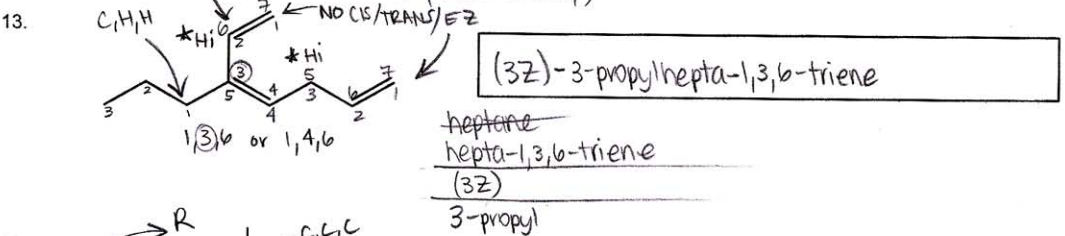
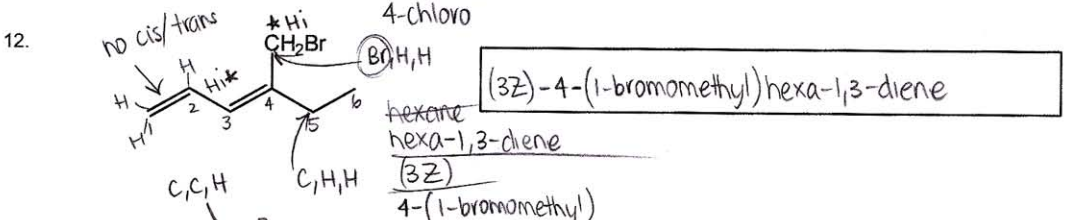
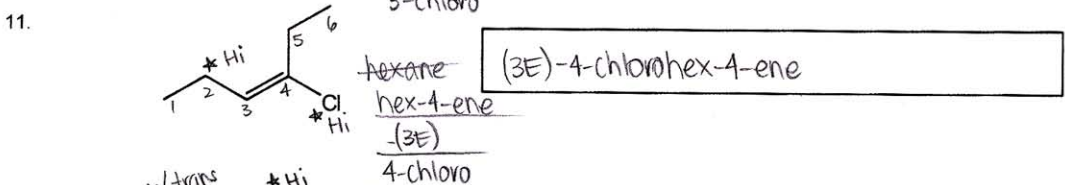
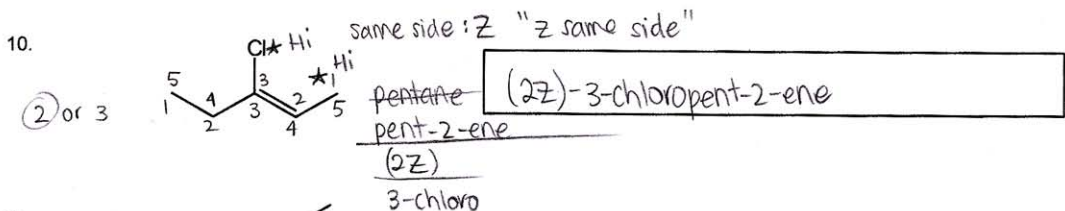
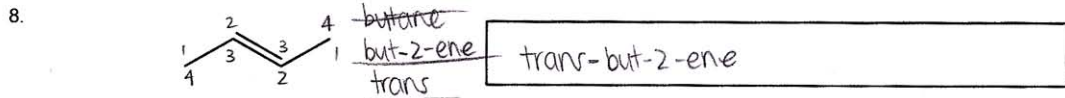


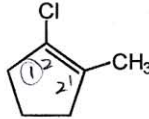
Chapter 7 Worksheet 1

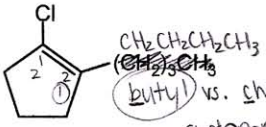
Give an acceptable IUPAC name for each of the following compounds. Be sure to indicate the **stereochemistry** where appropriate.

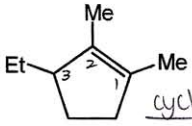


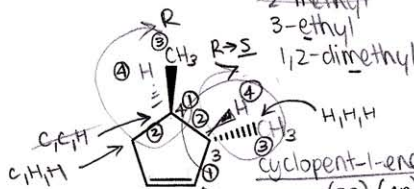


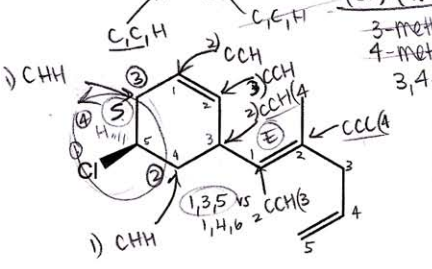
* can # from either side, resort to alphabet of substituents. "chloro" vs. methyl

15.  cyclopent-1-ene
1-chloro-2-methylcyclopent-1-ene

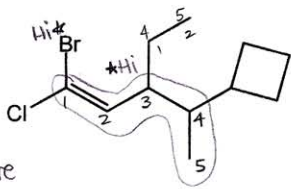
16.  cyclopent-1-ene
1-butyl-2-chloro

17.  cyclopent-1-ene
3-ethyl-1,2-dimethylcyclopent-1-ene

18.  cyclopent-1-ene
(3S, 4R)-3,4-dimethylcyclopent-1-ene

19.  cyclohex-1-ene
(5S)-5-chloro-3-((E)-2-methylpenta-1,4-dienyl)cyclohex-1-ene

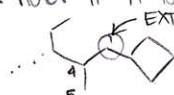
* The longest chain must contain @ least 1 double bond to even be considered as the parent chain.

20.  pent-1-ene
(1Z)-1-bromo-1-chloro-4-cyclobutyl-3-ethylpent-1-ene

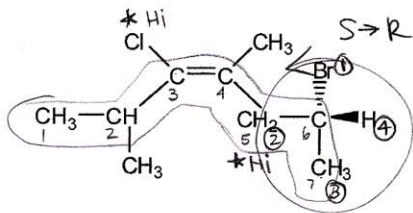
* The longest chain that contains as many double bonds as possible.

* # it to give the most substituents after finding longest chain because there are multiple "longest chains"

pent-1-ene
(1Z)
1-bromo
1-chloro
3-ethyl
4-cyclobutyl

* note: if it looked like this  that is not a simple cyclobutyl sub... that is a complex sub... 4-(1-cyclobutylmethyl)

21.



(3E, 6R)-6-bromo-3-chloro-2,4-dimethylhept-3-ene

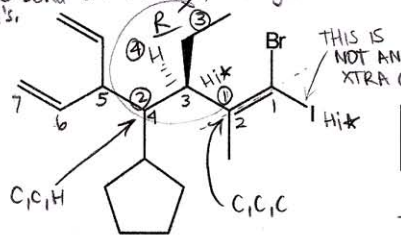
hept-3-ene

(3E) (6R)

- ~~2-methyl~~
- ~~3-chloro~~
- ~~4-methyl~~
- 6-bromo
- 2,4-dimethyl

★ Because you can # either way to get double bond on 1 and 6, # to give subs lowest #'s.

22.



(1E, 3R)-1-bromo-4-cyclopentyl-5-(1-ethenyl)-3-ethyl-2-methyl-1-iodohepta-1,6-diene

hepta-1,6-diene

(1E) (3R)

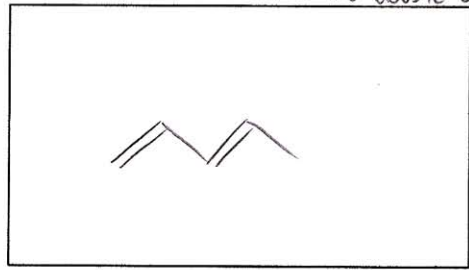
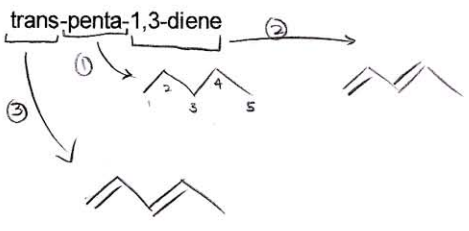
- ✓ 1-bromo
- ✓ 1-iodo
- ✓ 2-methyl
- ✓ 3-ethyl
- ✓ 4-cyclopentyl

✓ 5-(1-ethenyl)

★ was ethene, but it is a sub, so it is ethenyl w/ 1 indicating the pos of the double bond.

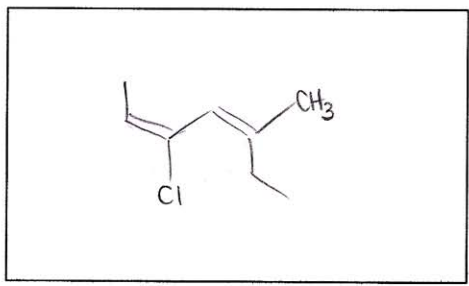
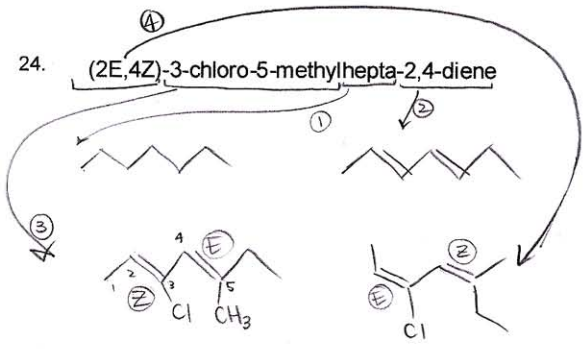
Draw the correct structure that corresponds to the following IUPAC names. Be sure to indicate **stereochemistry** where appropriate.

23.

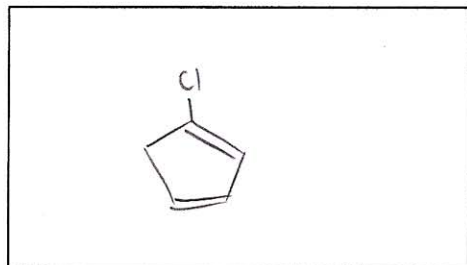
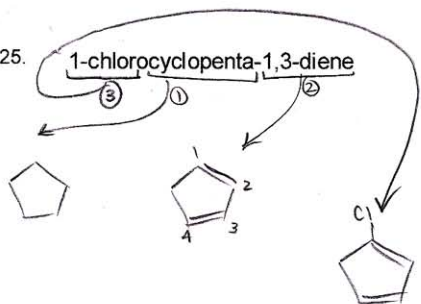


24.

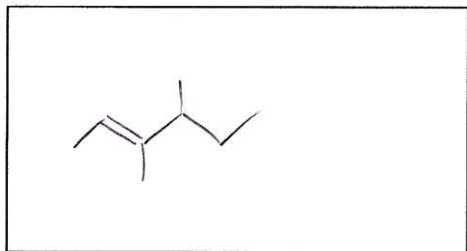
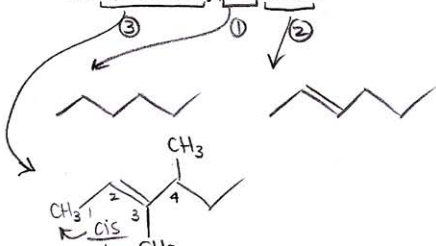
(2E, 4Z)-3-chloro-5-methylhepta-2,4-diene



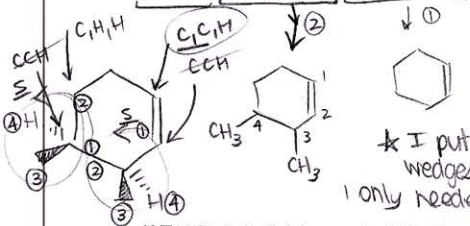
25. 1-chlorocyclopenta-1,3-diene



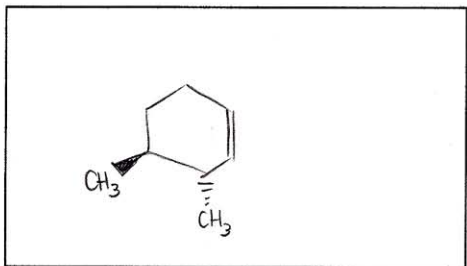
26. cis-3,4-dimethylhex-2-ene



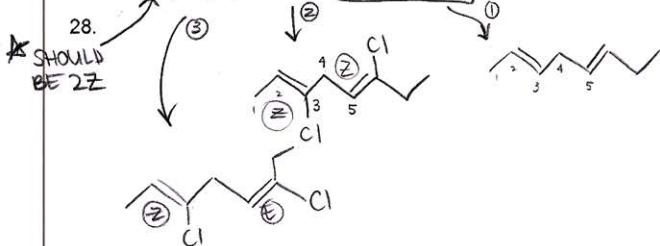
27. (3R,4S)-3,4-dimethylcyclohexene



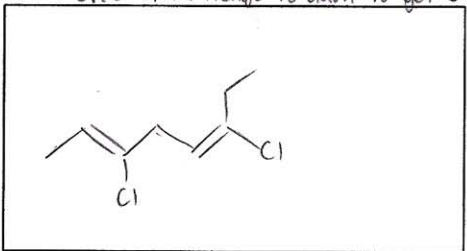
* I put them both on wedges initially + found I only needed to flip the CH₃ on carbon 3 from wedge to dash to get 3R.



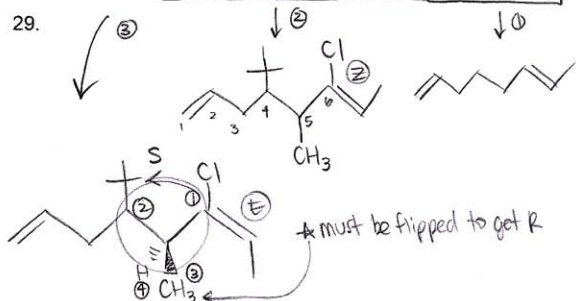
28. (3Z,5E)-3,6-dichloroocta-2,5-diene



* SHOULD BE ZZ



(5R,6E)-4-tert-butyl-6-chloro-5-methylocta-1,6-diene



* must be flipped to get R

