#### Beauchamp

On a separate piece of paper, provide acceptable names for the following structures. After you have named your structures, turn it around and redraw each structure by the name you gave it (for more practice). Clearly number the longest high priority chain in your drawing. Hints are provided just below the structures and answers are provided just below the hints. Use a separate piece of paper to cover those up unless you absolutely need them. There are probably some mistakes. Please let me know when you find them. Thanks. My usual 2 page overview of nomenclature is provided at the end of these problems (repeat found in other documents.)





Beauchamp









### 4. Problems - carboxylic acids



Beauchamp





# 5. Problems –anhydrides





Beauchamp









### 9. Problems - nitriles











15. Problems - miscellaneous low priority groups

#### Beauchamp

This problem has most of the functional groups you are responsible for in a first year course in organic chemistry. The problem, as difficult as it looks, is doable if you are systematic in your approach. A possible strategy to follow is listed just below.

## Some Points to Consider

- 1. What is the highest priority group present?
- 2. What is the longest chain with that group?
- 3. Number that chain (or ring) so that the high priority group gets the lowest number (often this is a 1 and not written).
- 4. Identify branches and lower priority functionality with their numbers (usually named as prefixes, except "-ene" and "-yne").
- 5. Combine everything into one name with proper use of hyphens (between numbers and letters) and commas (between numbers and numbers).



A possible solution (not alphabetized).



6-hydroxy-5-methylhexa-2E,4Z-dienyl (2S,6S,8Z,10R,11R)-2-carbamoyl-5-formyl-5-(3-heptyl-5,5-dimethylcyclopent-2-enyl)-6-mercapto-7,12-dioxo-8-nitro-9-butoxy-10-cyano-11-aminododec-8-en-3-ynoate



#### Beauchamp

Identify each carbon atom as primary, secondary, tertiary or quaternary.

Identify each carbon atom as methyl, methylene or methine.



Provide an acceptable structure for each of the following names.

a.	2,4-dimethylhexane							
b.	1,3,5,5-tetramethyl-1-cycloheptene							
c.	trans-5-hepten-1-yne	or		5E-hepten-1-yne				
d.	5-(4-methyl-2-cyclobutenyl)-1-nonene							
e.	trans-3-ethyl-1-isoproplycyclopentane							
f.	cis-1-t-butyl-4-isobutylcyclohexane							
g.	trans-2-cis-4-decadien-6,8-diyne		or	2E,4Z-decadien-6,8-diyne				
h.	1-allyl-3-propargyl-6-vinylcyclodecane	or	1-(2-proj	penyl)-3-(2-propynyl)-6-ethenylcyclodecane				
i.	1-hepten-6-yne							
j.	3-allyl-1-sec-butyl-1-cyclopropene	or	3-(2-propenyl)-1-(1-methylpropyl)-1-cycloprope					
k.	3-propyl-1-trans-6-octadiene		or	3-propyl-1E,6-octadiene				
1.	7-(1,1-dimethylethyl)-4-(1-methylpropyl)-1-cycloundecyne							
m.	2,9,10-trimethyl-6-butyldodecane							
n.	1-(1,3-dimethylbutyl)-4-methyl-1,3,5,7-cyclooctatetraene							
0.	2,9,10-trimethyl-6-butyldodecane							
p.	trans-(3,3-dimethylbutyl)-3-methylcyclooctane							
q.	7-(1,1-dimethylethyl)-4-(1-methylpropyl)-1-cycloundecyne							
r.	1-hepten-6-yne							
s.	1-trans-3-cis-7-cis-cycloundecatriene	or		1E,3Z,7Z-cycloundecatriene				
t.	3-benzyl-6-ethyl-7-phenyl-2Z-octene							

Beauchamp

## Some aromatic structures and their common names (not required in 314 or 315).



# **General Strategy For Naming Simple Organic Compounds (Bare bones summary sheets)**

- 1. Find the highest priority group. These are listed in order of priority in the table of functional groups.
- 2. Find the longest chain containing the highest priority group. You should know carbon chains of length  $C_1$ - $C_{12}$  (listed in the table).
- 3. Number the longest chain containing the highest priority group to give the highest priority group the lowest number possible in numbering the longest chain. For the first seven groups, the functional group carbon will be number 1 (if it is the highest priority group) and the "1" can be omitted, since it is understood that it has to be this way.
- 4. Usually the highest priority group is named as a suffix at the end of alkane, alk-#-ene or alk-#-yne. The final e is dropped if the suffix begins with a vowel and it is retained if the suffix begins with a consonant. A number will be present in front of the suffix name unless its position is unambiguously clear (i.e. carboxyl groups aldehydes, nitriles, etc. always = 1, if highest in priority). If there is a C/C pi functional group to identify (alkene or alkyne), the number in front of its part of the name describes its position (see rule 6 below). If both a pi bond and a high priority substituent are present, then two numbers may be necessary, one for each functionality.
- 5. Lower priority groups are named with their prefix names and their location numbers based on the numbering of the parent chain (always true for substituents numbered 12 on the next page). The lower priority substituents should be listed in alphabetical order. Some parts of prefix names count in this regard and some don't. We will not emphasize this aspect in this course.
- 6. Double bonds and triple bonds are named as alk-#-ene or alk-#-yne, respectively. If both are present, name as alk-#-en-#-yne. Multiple pi bonds (or other substituents) use the prefixes di, tri, tetra, penta etc. with a number for each occurrence. In such cases, an "a" is added in front of the numerical prefix for better phonetics. (alka-#,#-diene or alka-#,#,#-triyne, alka-#,#-diene, etc.)

The essential functional groups to know (for our course) and their prefixes and suffixes are given in the attached table. In this table the term "alkan-#-suffix" is a generic term for any alkane with a functional group suffix, and it must be replaced with the correct parent stem name based on the number of carbons in the longest chain ( $C_1$ - $C_{12}$  for us). If there is a double bond, the name will change to "alk-#-en-#-suffix and if there is a triple bond, the name will change to "alk-#-en-#-suffix.





(2Z,4S)-3-ethoxy-4-methyl-8-oxooct-2-en-6-ynoic acid

Notice that no "1" is used for the carboxylic acid group, because it has to be "1" in this structure.

Beauchamp

<b>Functional Group</b>	0	<u>prefix</u>	The part of each name specific to the functional							
1. Carboxylic acid		#-carboxy not considered	alkan <u>oic acid</u>	group is in <b>bold</b> and <u>underlined</u> to help you see those features						
2. Anhydride		#-acyloxyalkanecarbonyl not considered	alkan <u>oic anhydride</u> (if symmetrical)	"R" = carbon chains						
3. Ester	к 0 к 0 " " "	#- <u>alkoxycarbonyl</u>	<u>alkyl</u> alkan <u>oate</u> (R') (RCO <sub>2</sub> )	# carbons	alkane chain name	alkyl branch name				
4. Acid halide	R <sup>C</sup> CI	#- <u>chlorocarbonyl</u>	alkan <mark>oyl chloride</mark>	1 2 2	methane ethane	methyl ethyl				
5. Amide	0 " R <sup>-</sup> NH <sub>2</sub>	#- <u>carbamoyl</u> # <u>-amido</u>	alkan <b>amide</b>	5 4 5 6	butane pentane bexane	propyi butyl pentyl bexyl				
6. Nitrile	R−C≡N	#- <u>cyano</u>	alkane <u>nitrile</u>	7 8 9	heptane octane	heptyl octyl nonvl				
7. Aldehyde	о "С к^ <sup>С</sup> `Н	#- <u>oxo</u>	alkan <u>al</u>	10 11 12	decane undecane	decyl undecyl dodecyl				
8. Ketone	0 " R <sup>^C</sup> R'	#- <u>oxo</u> (older = #- <u>keto</u> )	#-alkan <u>one</u>	13 14 15	tridecane tetradecane	tridecyl tetradecyl pentadecyl				
9. Alcohol	R-OH	#-hydroxy	#-alkan <u>ol</u>	16 17 18	hexadecane heptadecane	hexadecyl heptadecyl				
10. Thiol	R-SH	#- <u>mercapto</u>	#-alkane <mark>thiol</mark>	19 20	nonadecane icosane etc.	nonadecyl				
11. Amine	R-NH <sub>2</sub>	#- <u>amino</u>	#-alkyl <u>amine</u> /#-alka	an <b>amine</b>						
12. Ether R-O-R'		#-alk <u>oxy</u> (if more than 5C's, then #-alkoxyl) (can also use "#-oxa" prefix and								
12. Thioether	R-S-R'	count as carbo #-alkyl <u>thio</u>	n in longest chain)							
12. Halogen	R-X	#- <u>fluoro</u> , #- <u>chloro</u> , #- <u>bro</u>	> always prefixes (no suffix names)							
12. Azide*	R-N <sub>3</sub>	#- <u>azido</u>								
12. Diazo*	R-N <sub>2</sub>	#- <u>diazo</u>								
12. Nitro*	R-NO <sub>2</sub>	#- <u>nitro</u>								
12. Nitroso	R-NO	#- <u>nitroso</u>								
12. Carbon branches R— #-alkyl, #-(alk-#-enyl), #-(alk-#-ynyl) * = formal charge is necessary in these Lewis structures and there are two reasonable resonance structures										
A A A A A A A A A A A A A A A A A A A										
riority functional above -yne drop preceeding "e" if groups with location # with location # suffix starts with a vowel										

Z:\classes\314\314 Special Handouts\Nomenclature programed probs.doc