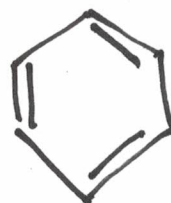
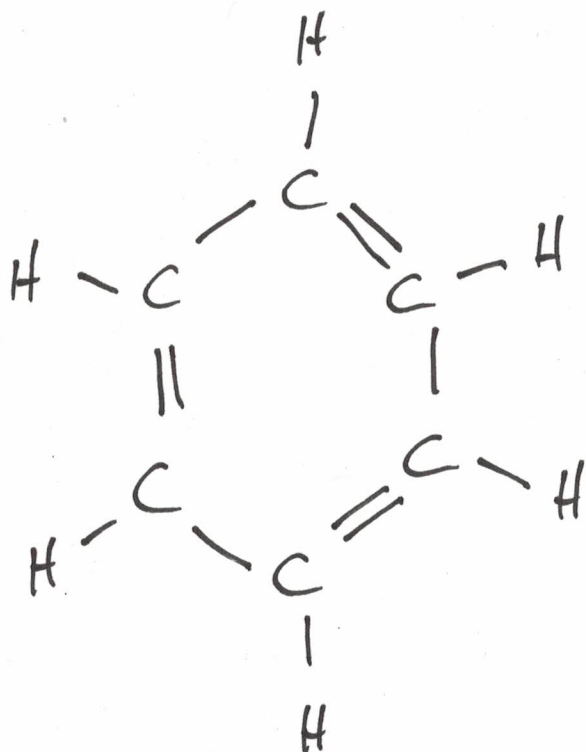
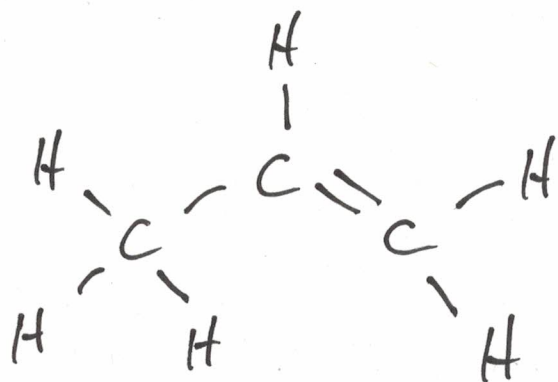
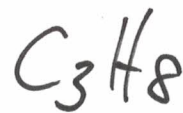
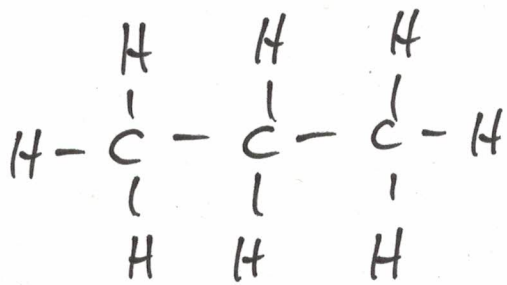
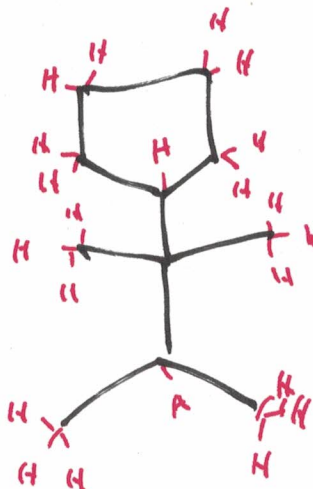
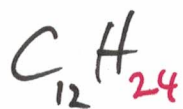
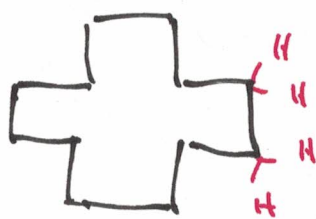
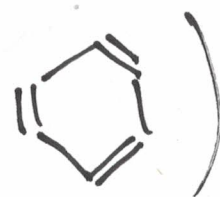


Skelett / Summenformel / (e-)




Cardo i.o. :



- * 1 Nachbar ...
- 2 Nachbarn → 60°
- 3 " → 120°
- 4 " → 110° (gezeichnet 90°)

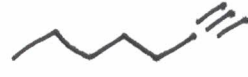
Bindungsmöglichkeiten

C-C; Einfachbindung, z.B. 

Alkan









C=C; Doppelbindung, z.B. 

Alken

C≡C; Dreifachbindung, z.B. 

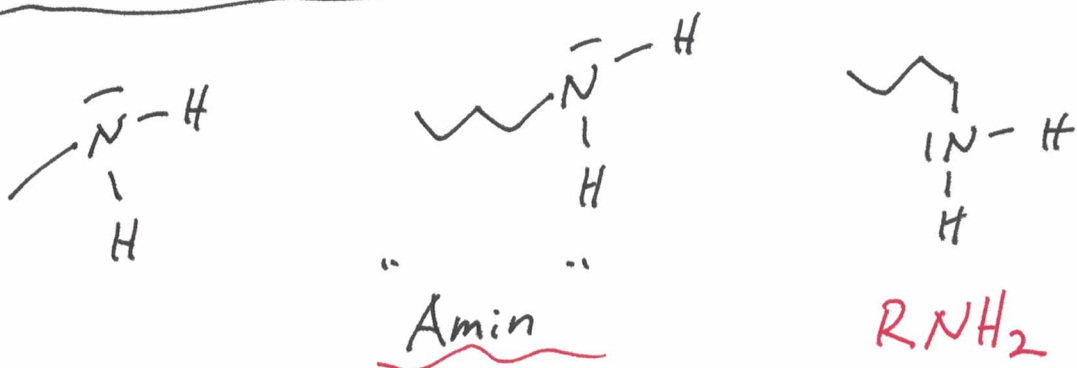
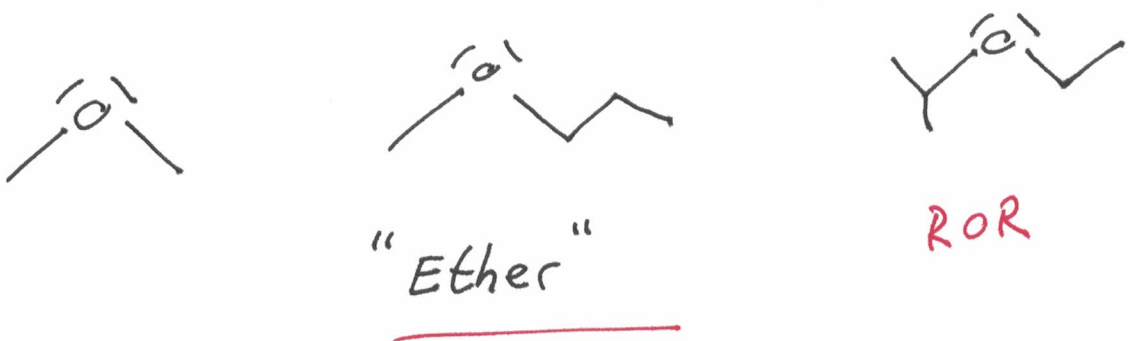
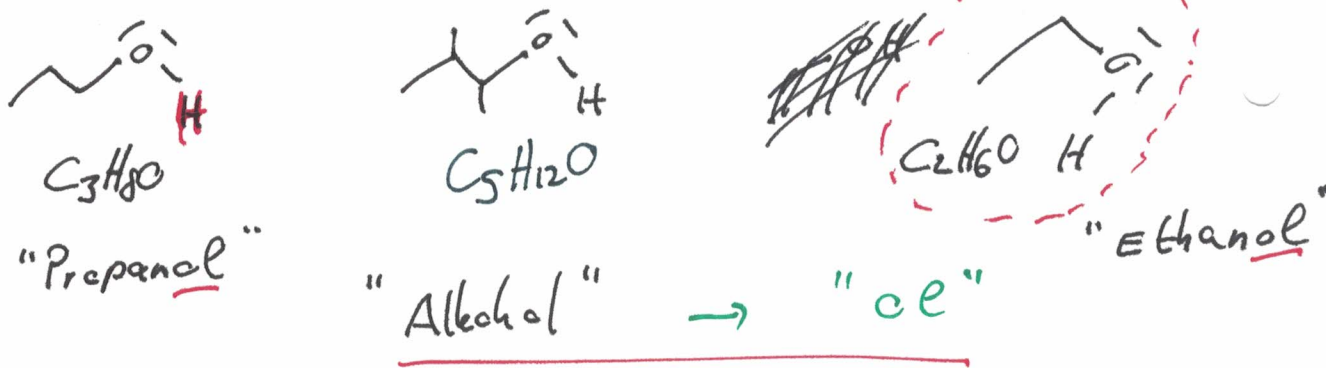
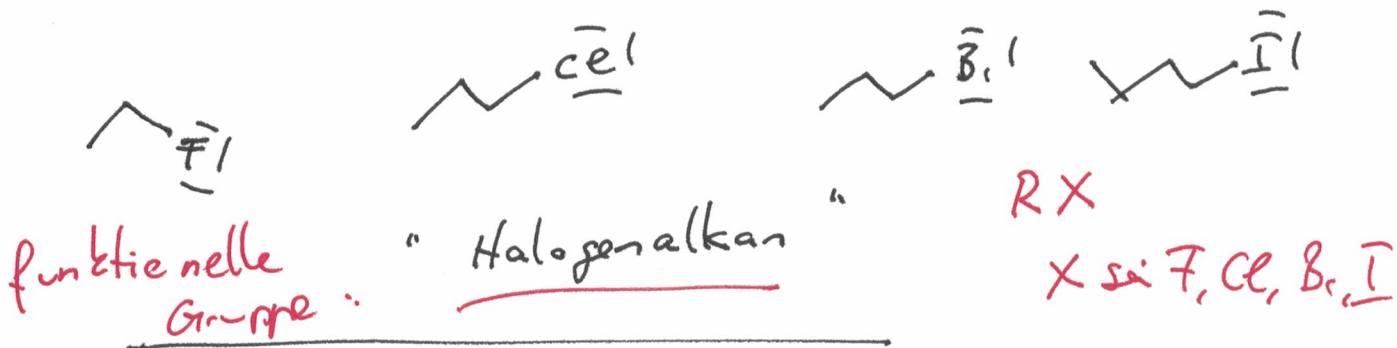
Alkin

Anzahl C-Atome

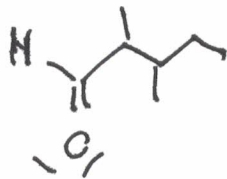
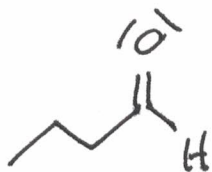
gas- fö- rig	1	$\begin{array}{c} \text{H} \\ \\ \text{H}-\text{C}-\text{H} \\ \\ \text{H} \end{array}$	Methan	2	/	Ethan
	3		Propan	4		Butan
	5		Pentan	6		Hexan
flüssig	7		Heptan	8		Octan
	9		Nonan	10		Decan

fest | ~ 20 - C-Atome

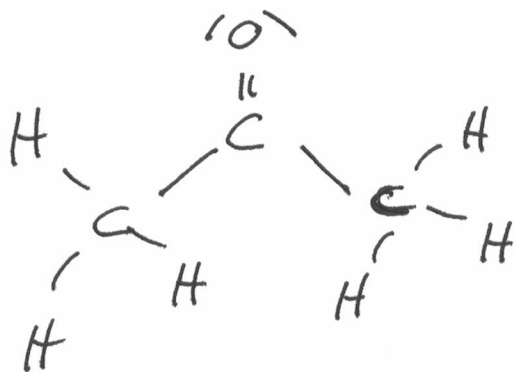
• " Eine Einfachbindung zwischen C und einem elektronegativeren Atom "



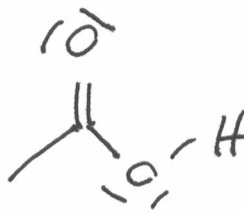
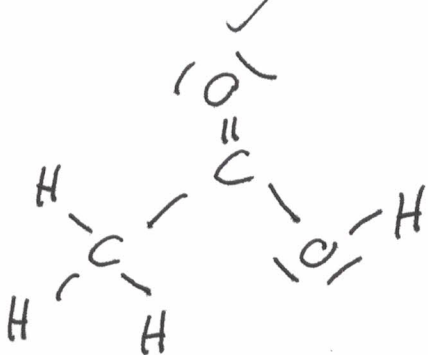
• Doppelbindung zwischen C + O: $C=O$



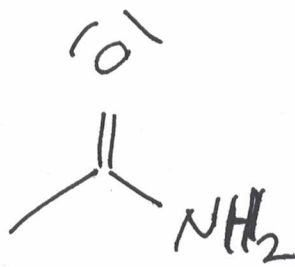
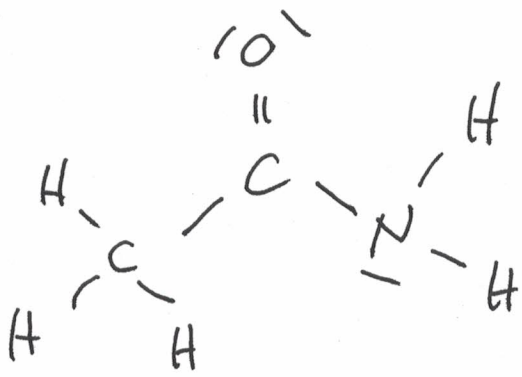
Aldehyd



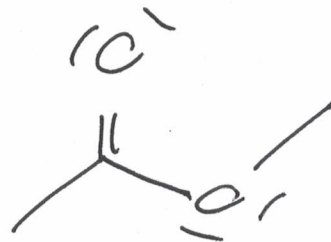
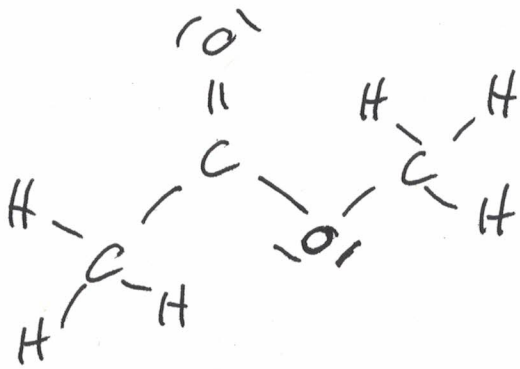
Keton



Carbonsäure



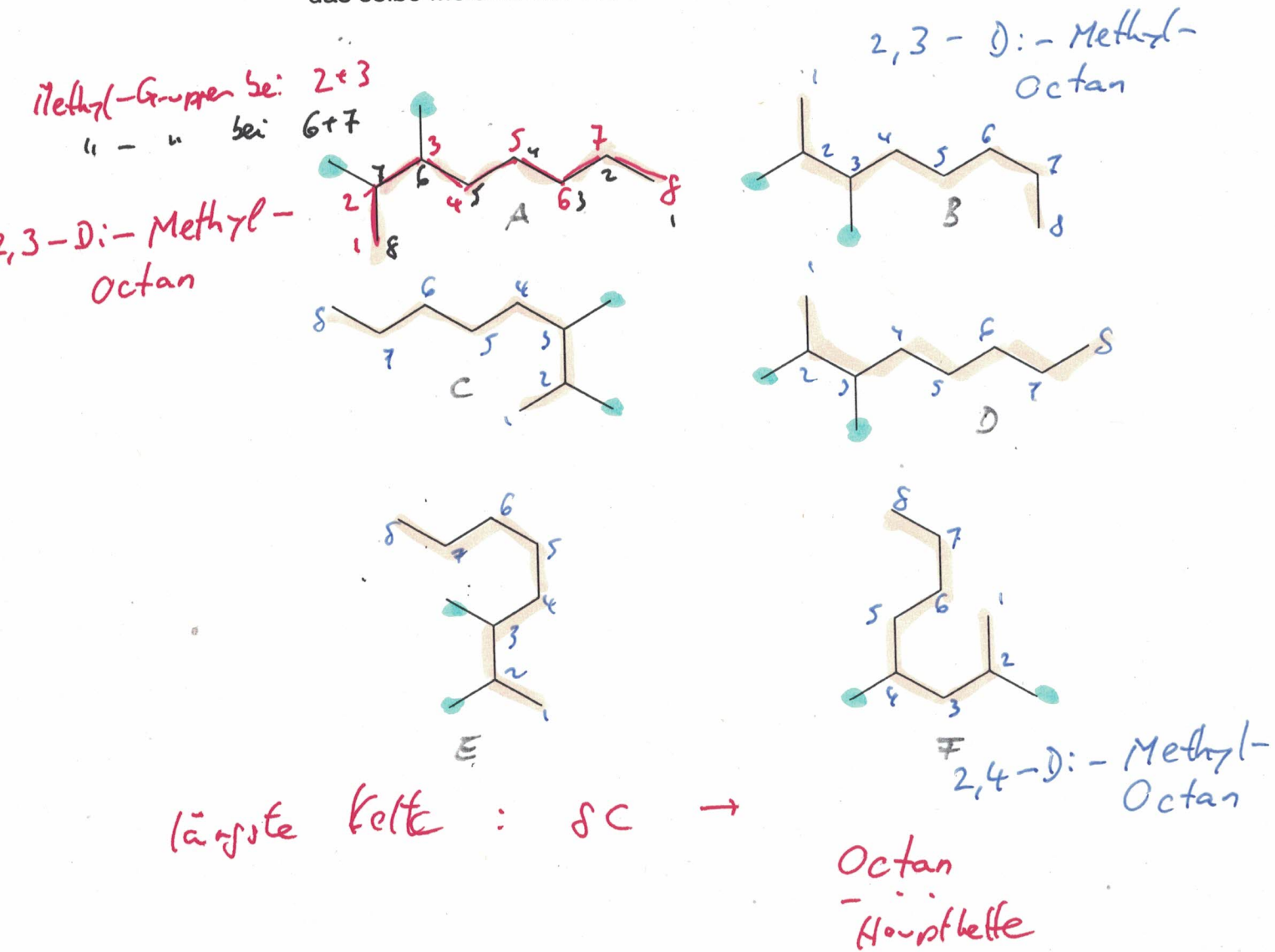
"Amid"

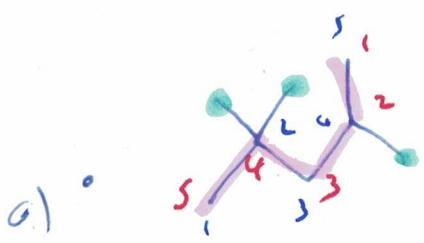


Ester



2.1 Gemäss der Theorie der kovalenten Bindungen herrscht um jede C-C-Einfachbindung freie Drehbarkeit. Kohlenwasserstoff-Moleküle sind von daher sehr beweglich. Logischerweise kann man ein bestimmtes Kohlenwasserstoff-Molekül auf sehr viele Arten aufzeichnen. Die folgenden Skelettformeln stellen bis auf eine Ausnahme alle das selbe Molekül dar. Welches Molekül ist hier fehl am Platz?





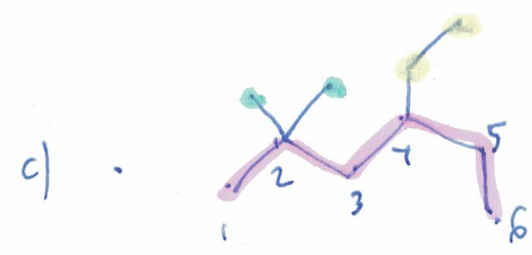
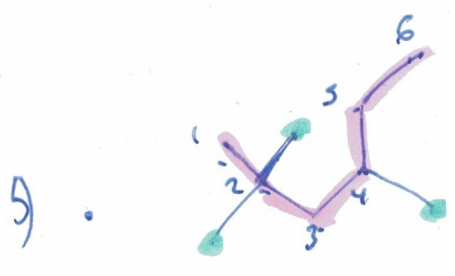
• 5C Pentan

• 3x CH₃-Rest

2,4,4-Tri-Methyl-Pentan
2,2,4 "

• 6C Hexan

• 3x CH₃-Rest



• 6C Hexan

• 2x CH₃-Rest

• 1x CH₂-CH₃

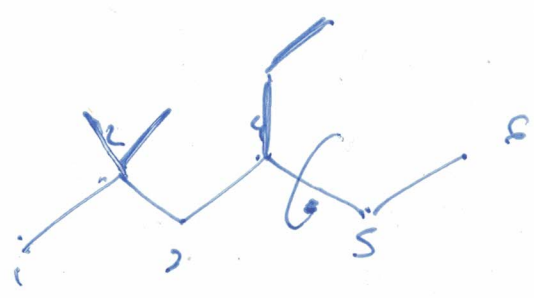
Methyl

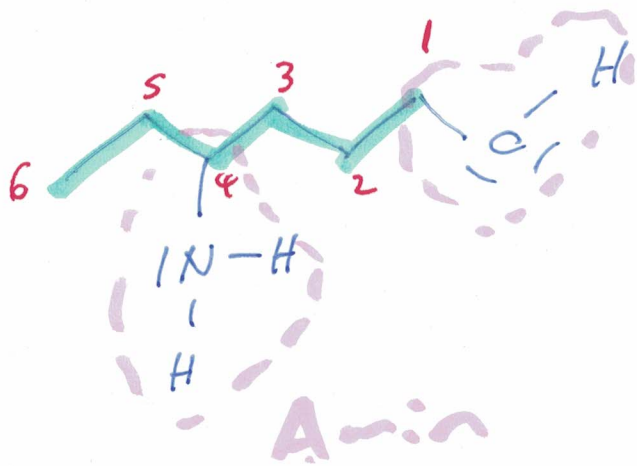
Ethyl

a) 2,2,4-Tri-Methyl-Pentan

b) 2,2,4 Tri-Methyl-Hexan

c) 4-Ethyl - 2,2-Di-Methyl - Hexan





Alkohol

• 6C → Hexan

Position 1 : Alkohol

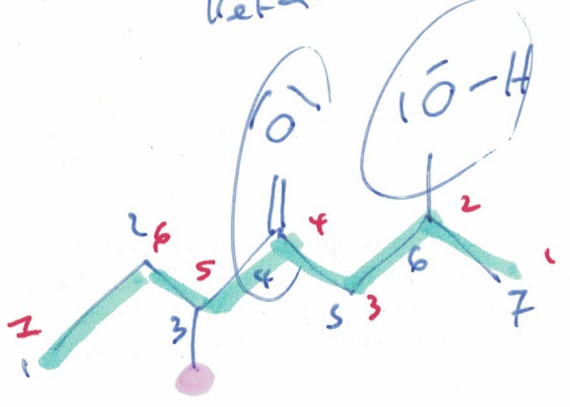
„ 4 : Amin

„ 4-Amino - Hexan - 1-ol „

„ Amin - Hexan - Alkohol

Präfix - "Hauptkette" - Suffix

	Präfix	Suffix
fun. Gruppe		
Alkohol	Hydroxy -	-ol
Keton	Oxo -	-on
Aldehyd	Formyl -	-al
Amin	Amino -	-amin



AlE

3-ethyl-

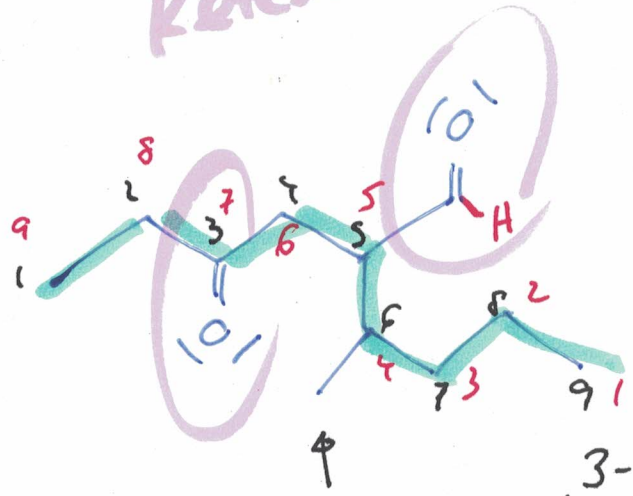
6 Hydroxy - Heptan - 4-on

Methyl - Rest

2-Hydroxy - 5-Ethyl - Heptan - 4-on

Keten

Aldehyd!



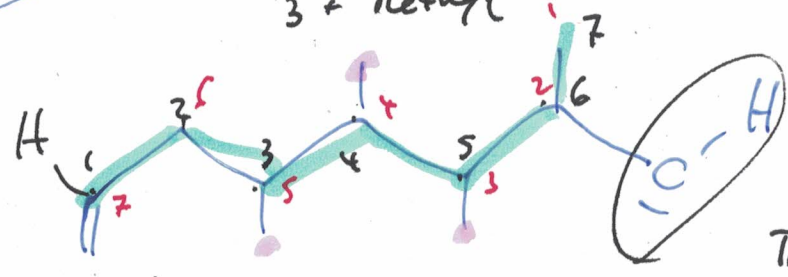
9 C - Nonan

Oxo - Nonan - al

Ethyl

3-oxo - 6-ethyl - Nonan - 5-al
(7-Carb - 4-Ethyl - Nonan - 5-al)

3 + Ethyl

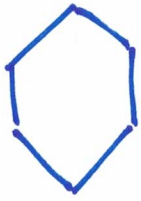


Aldehyd!

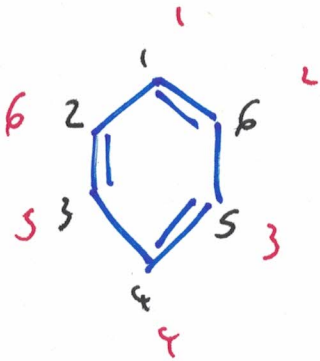
~~6 C - Hexan~~
7 C - Heptan

Trimethyl

6-Hydroxy-2,3,5-Trimethyl-Heptan-1-al
(2-Hydroxy-3,4,5-Trimethyl-Heptan-7-al)



Cyclo - Hexan

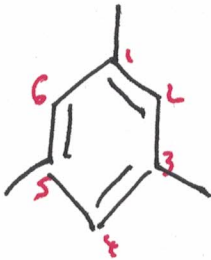


Cyclo - Hexen

Cyclo - Hex - Tri - en

Cyclo - Hex(an) - Tri - 2,4,6 - en

Cyclo - Hex(an) - Tri - 1,3,5 - en

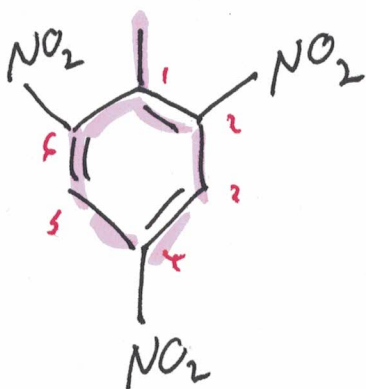


1,3,5 - Tri - ethyl -

cyclo - Hexan - Tri - 1,3,5 - en

TNT

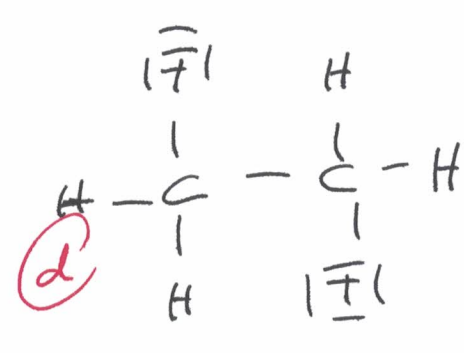
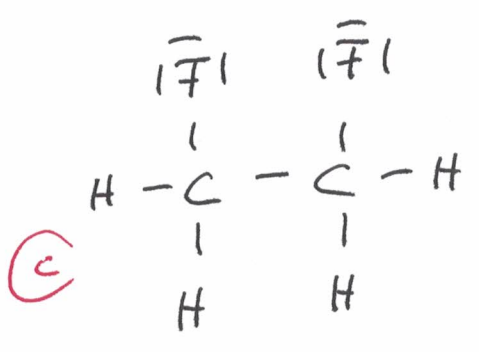
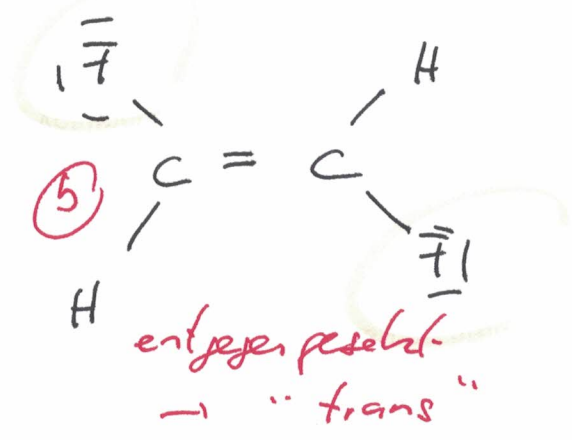
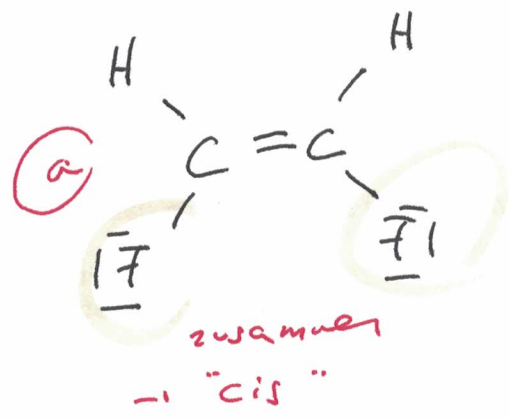
Tri - Nitro - Tolvol



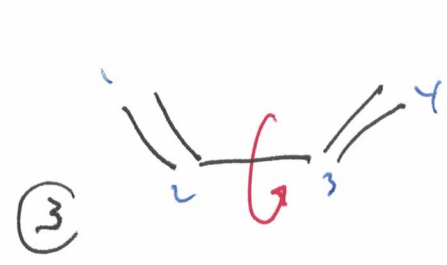
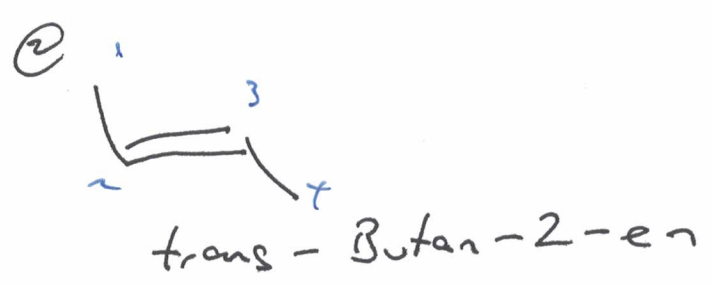
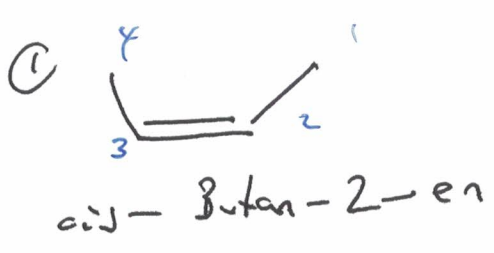
2,4,6 - Tri - Nitro - 1 - ethyl -

cyclo - Hexan - Tri, 1,3,5 - en

Bsp.



c ≡ d



3 ≡ 4 (!)

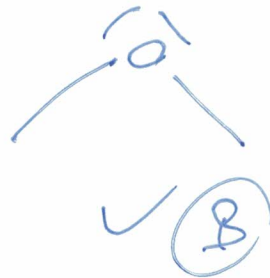
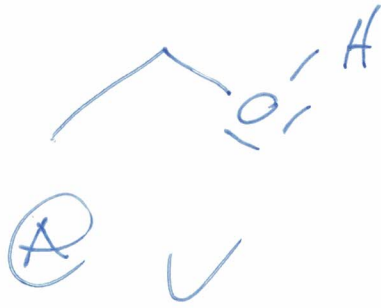


Butan-1,3-Di-en

verschiedenes räumliches Aufbau,
ABER.. gleiche Summenformel!

→ "Isomere"

• C_2H_4O

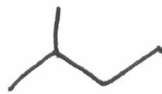


gleiche Summenformel,
verschieden funktionelle

Gruppen →

Funktions-
isomere

• C_5H_{12}



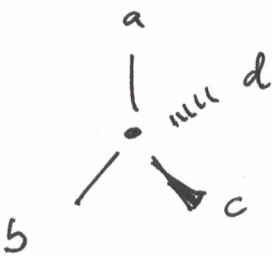
gleiche Summenformel,
unterschiedliches Skelett

→ Skelett-isomere

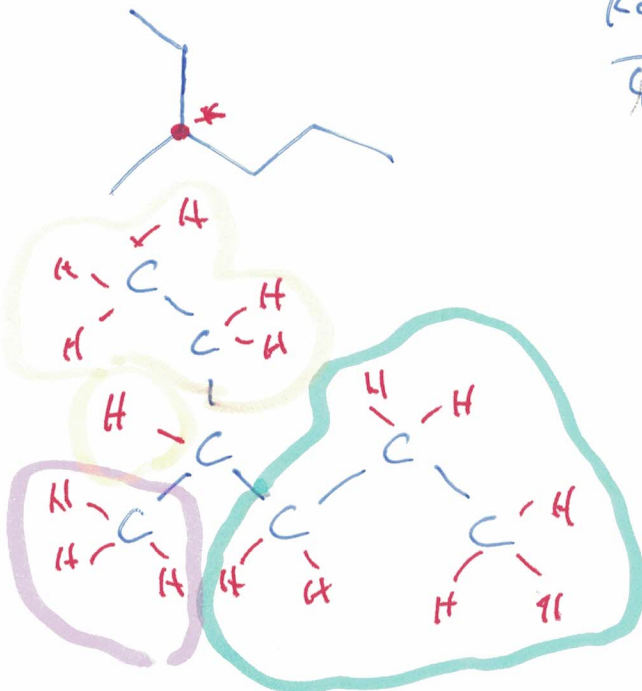
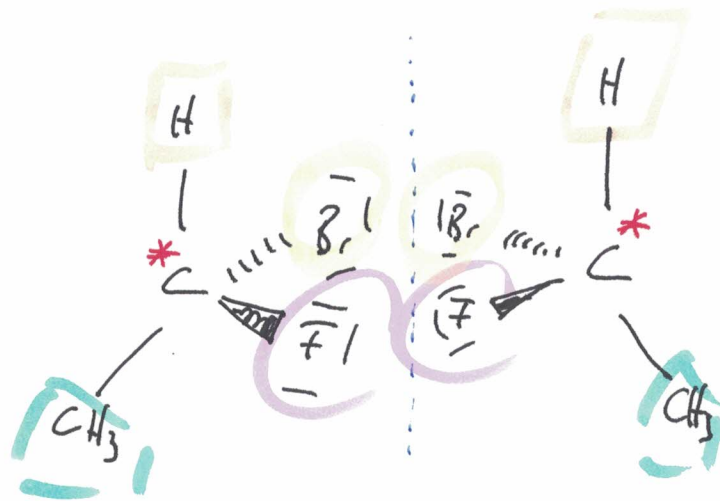
Chiralität

"Etp."

Die Chiralität nennt man die Eigenschaft bestimmte Gegenstände (Moleküle), deren Bild resp. Spiegelbild durch Drehung sich nicht zur Deckung bringen lassen



z.B.



Kohlenstoffatom(!) mit
4 verschiedenen Resten?

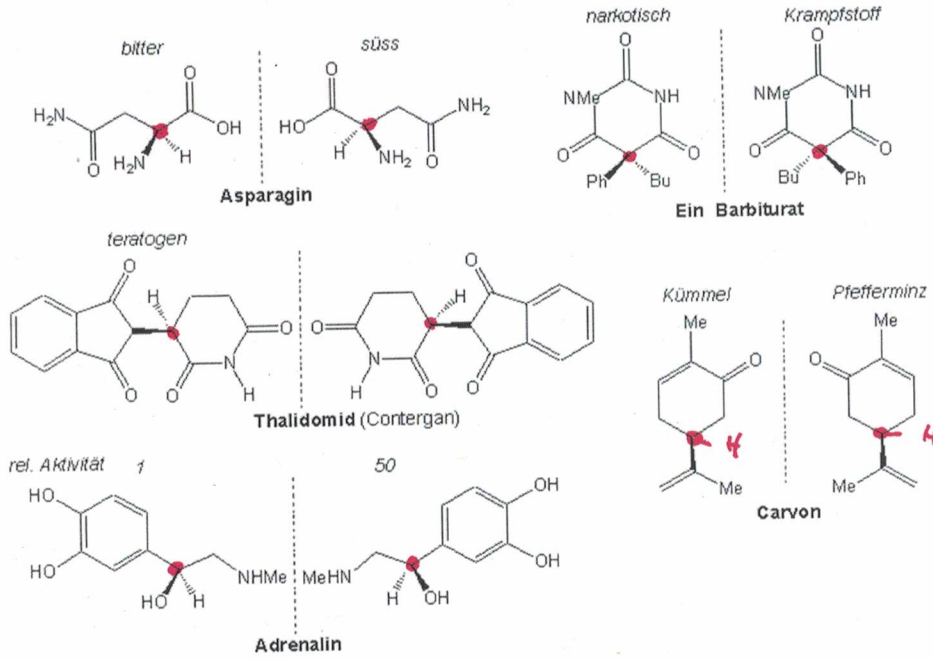
ja: Chiralitätszentrum *

Chiralität

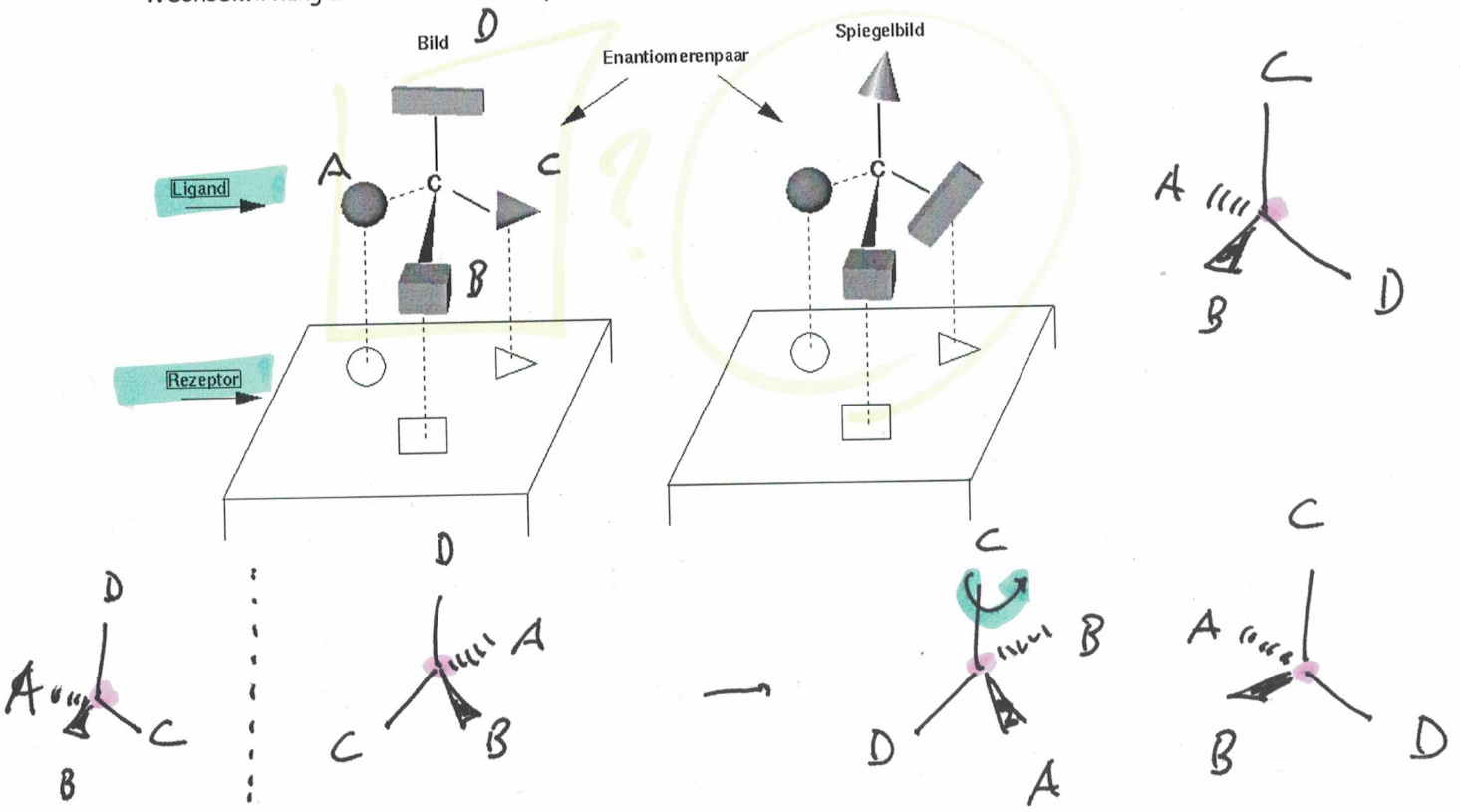
Chiralität in der Natur:

Betrachtet man Moleküle in der freien Natur, so wird sehr schnell ersichtlich, dass praktisch alle wichtigen Substanzen chiral sind. Beispiele:

• Chiralitätszentrum

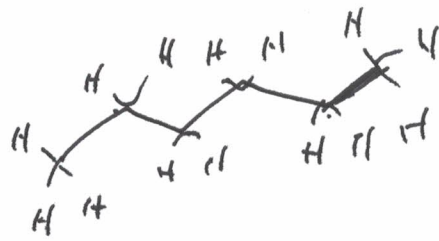


Warum haben solche Stereoisomere unterschiedliche biologische Eigenschaften? Um ihren biologischen Effekt auszuüben müssen die Moleküle mit anderen Biomoleküle (Proteinen oder DNA, z.B.) Komplexe bilden. Die Biomoleküle sind chiral! Ein chirales Molekül muss zu seinem spezifischen Rezeptor passen (wie eine Hand zu einem Handschuhe!). Zum Beispiel können wir in der folgenden Abbildung eine spezifische Wechselwirkung zwischen dem Rezeptor für Adrenalin und den zwei Adrenalin Stereoisomeren darstellen:

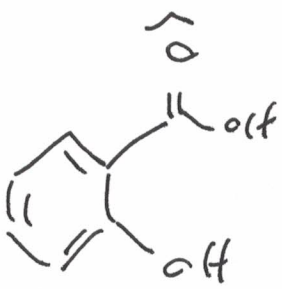


Spiegel-
bild

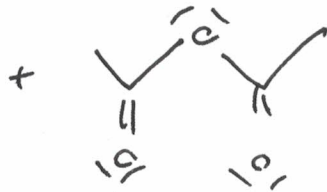
C_6H_6



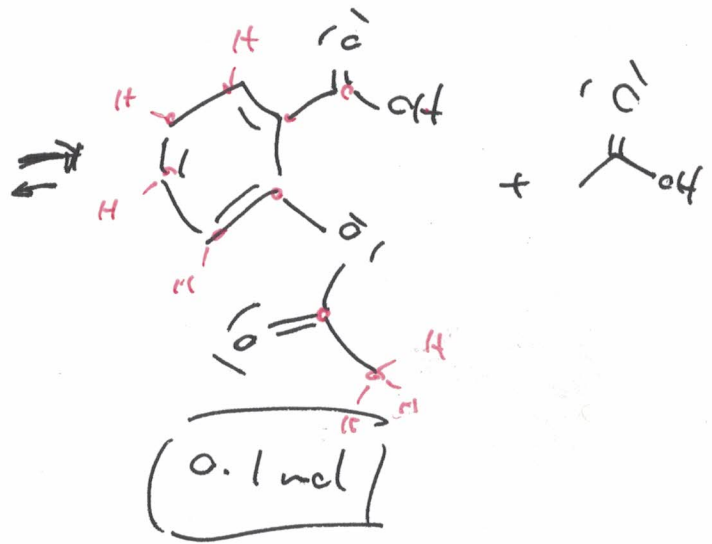
↳ Lösung:
zeigen.



0.1 mol



0.12 mol



Succinylsäure



$$M(C_9H_8O_4) = 9 \cdot 12 + 8 \cdot 1 + 4 \cdot 16 = 180 \text{ g/mol}$$

$$1 \text{ mol} \hat{=} 180 \text{ g}$$

$$0.1 \text{ mol} \hat{=} \underline{\underline{18 \text{ g}}}$$

maximale Ausbeute 100%

Praxis : ~ 90% Ausbeute



+

