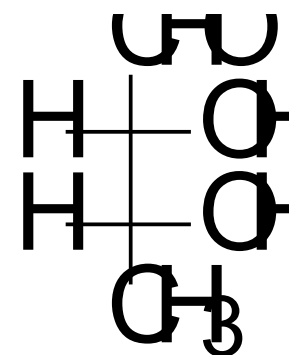
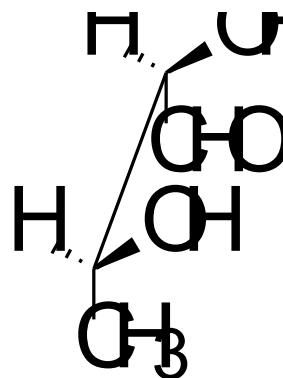
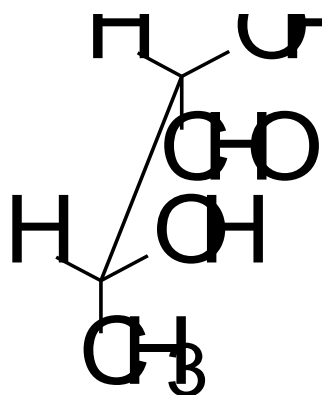
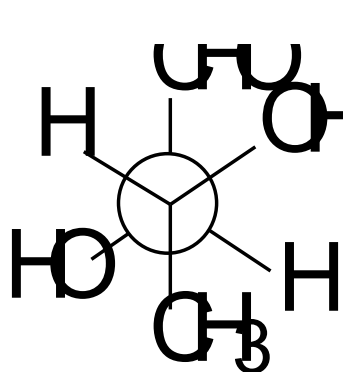


# STEREOCHEMIE

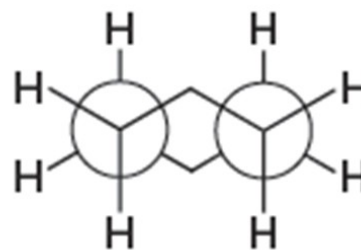
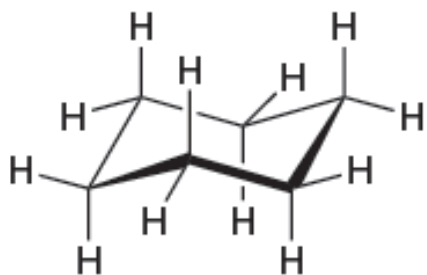
# Typy vzorců

Následující vzorec v Newmanově projekci přepište do Fischerovy projekce



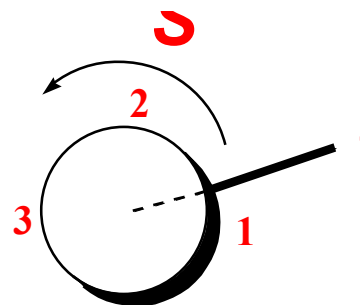
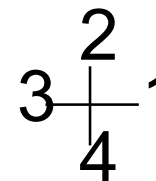
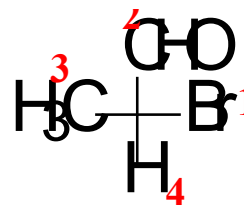
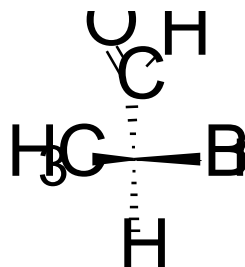
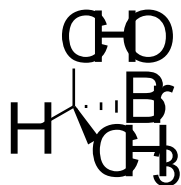
# Typy vzorců

Molekulu cyklohexanu překreslete do Newmanovy projekce

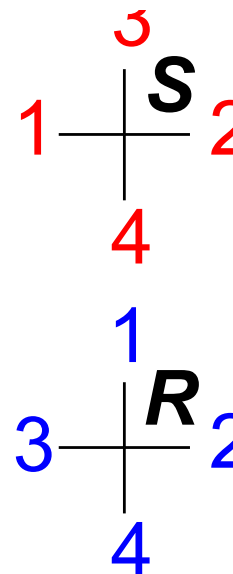
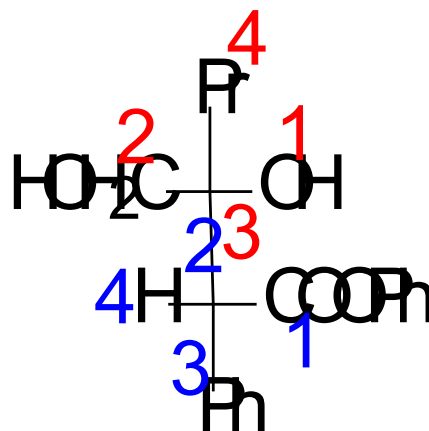
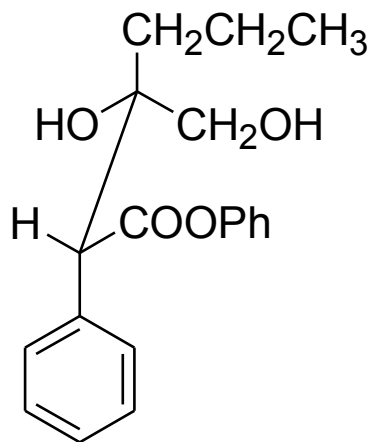


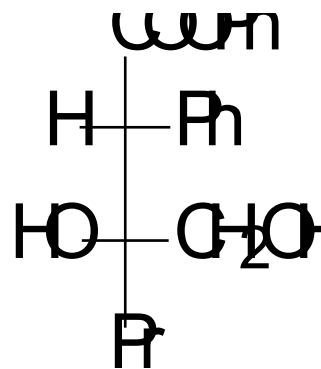
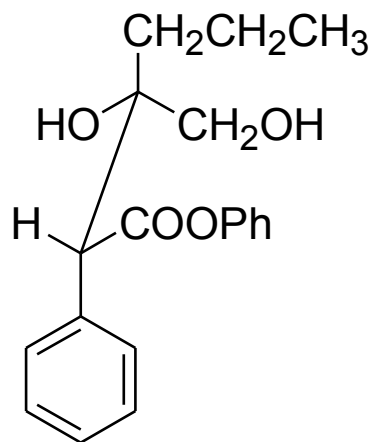
# Vytvořte systematický název uvedené sloučeniny

(2S)-2-bromopropanal



U následující sloučeniny identifikujte centra chiralidy a určete na nich absolutní konfiguraci

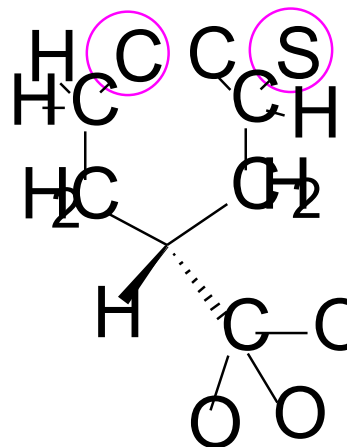
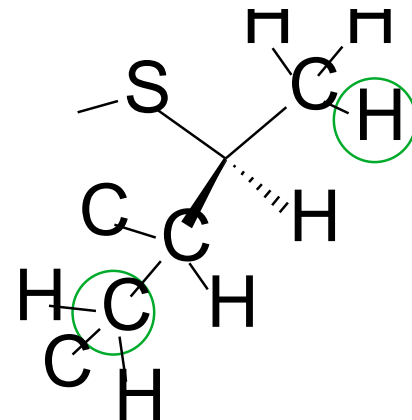
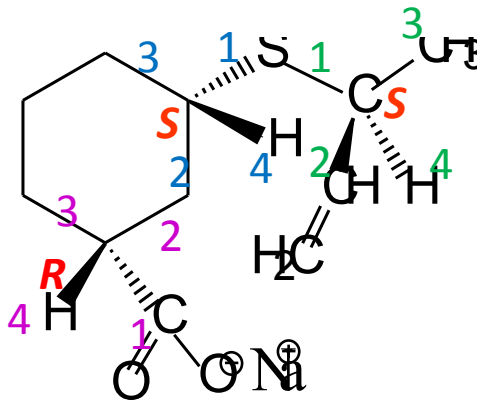
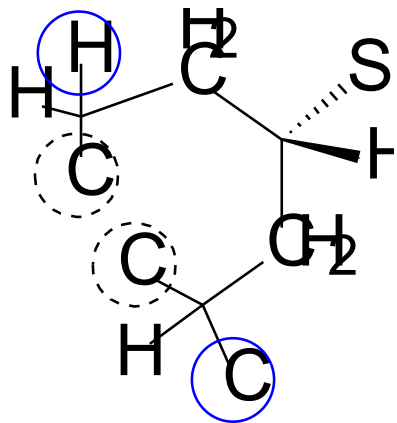




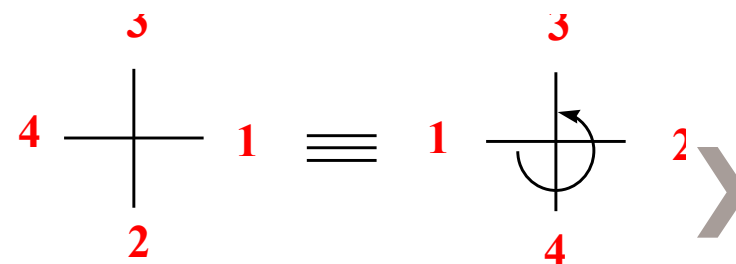
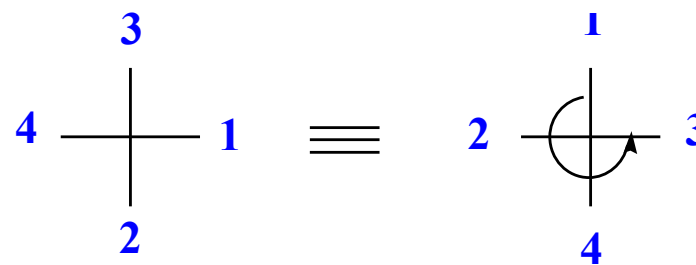
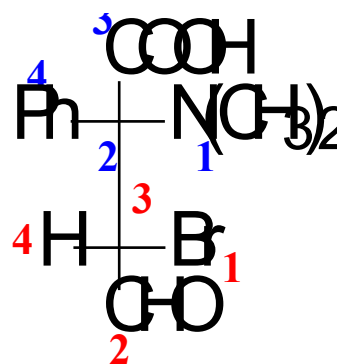
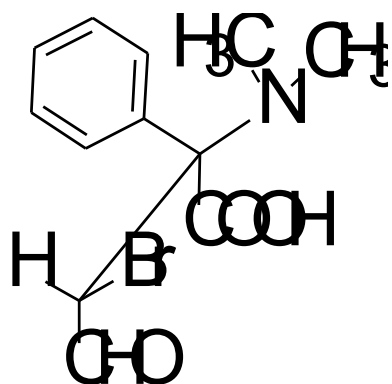
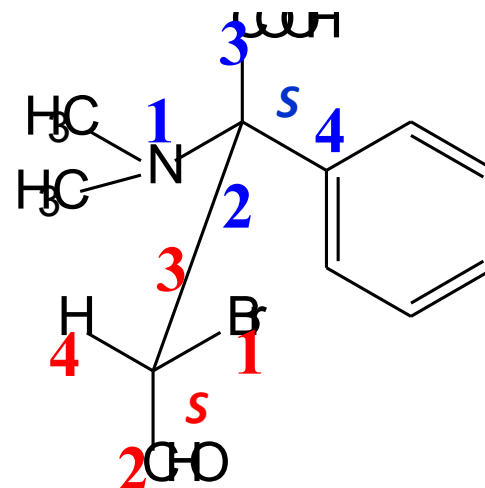
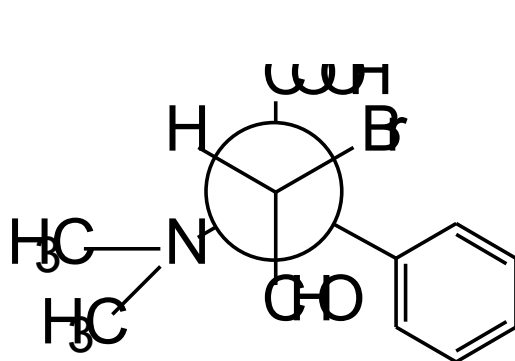
Správně zakreslená Fischerova projekce



# Označte centra chiraloty a určete na nich absolutní konfiguraci



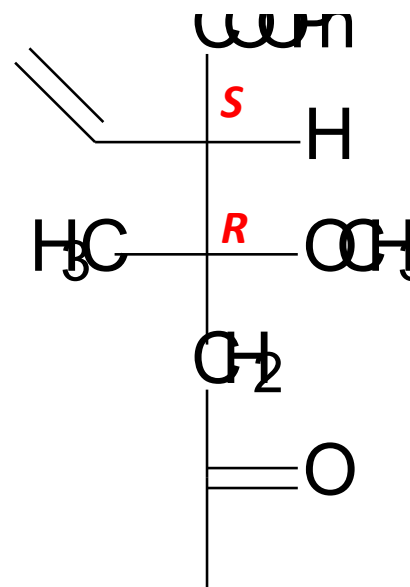
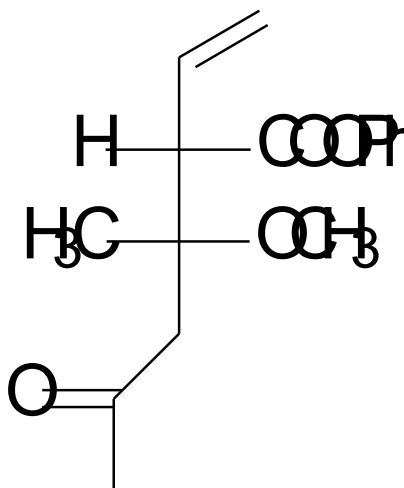
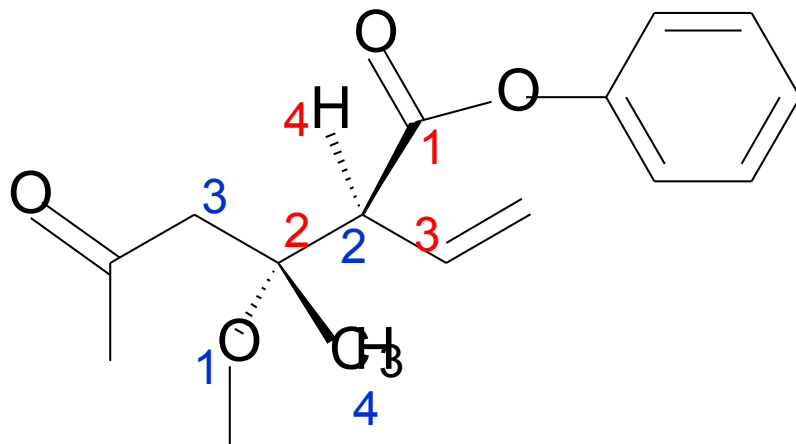
Najděte v molekule centra chirality a určete na nich absolutní konfiguraci a sloučeninu pojmenujte



(2S,3S)-3-brom-2-(N,N-dimethylamino)-2-fenyl-4-oxobutanová kyselina



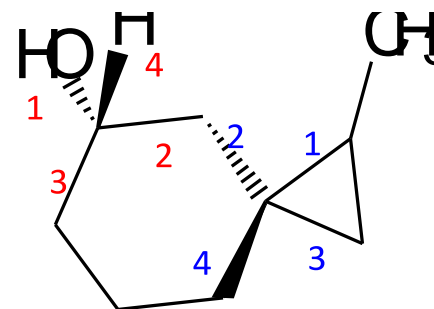
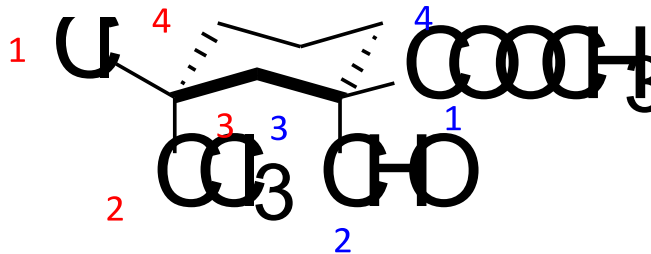
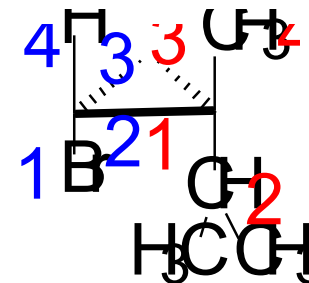
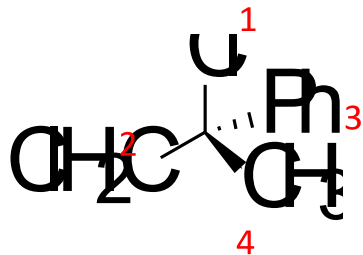
Molekulu překreslete do Fischerovy projekce a určete absolutní konfiguraci na centrech chiralidy



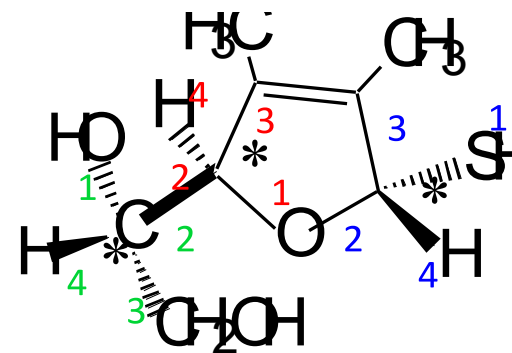
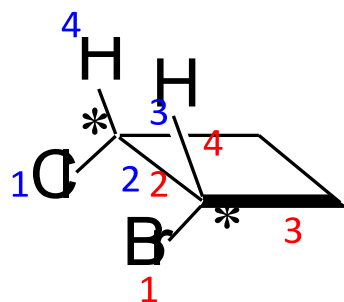
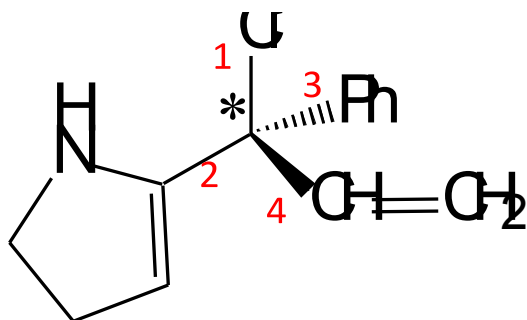
(2*S*,3*R*)-fenyl-3-methoxy-3-methyl-5-oxo-2-vinylhexanoát



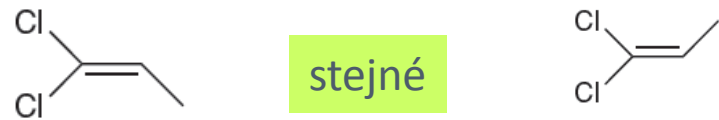
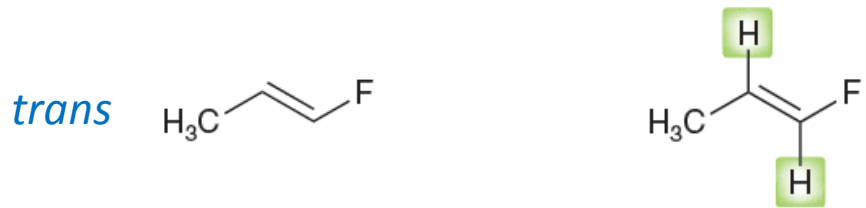
Označte \* centra chiralidy a určete na nich absolutní konfiguraci



Na vyznačených centrech chirality určete absolutní konfiguraci



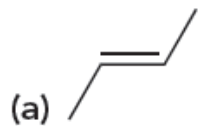
# » Geometrické izomery



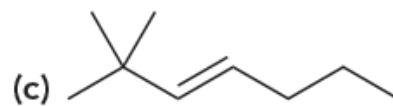
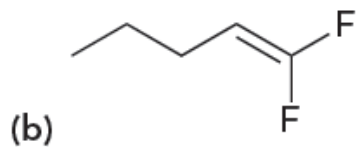
stejné



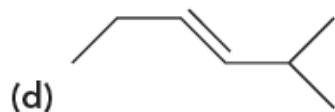
Kde je to potřeba, určete, o jaký stereoizomer se jedná (*cis - trans, E - Z*)



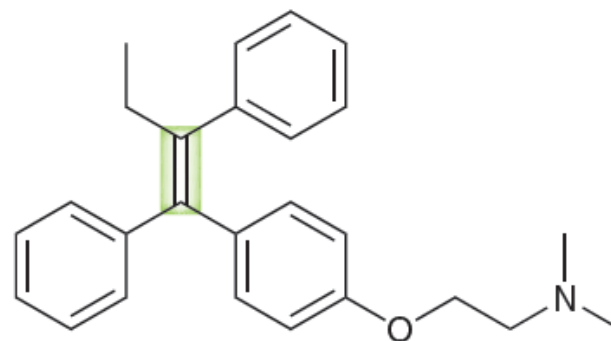
*trans*



*trans (E)*



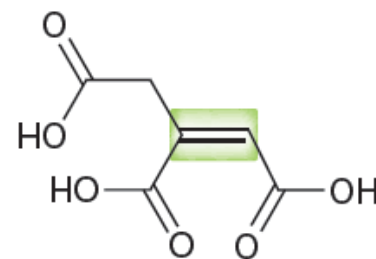
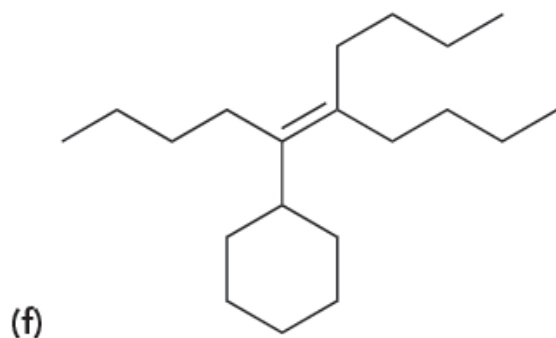
*trans (E)*



**Tamoxifen**

Used in treatment of breast cancer

*trans (Z)*



**Aconitic acid**

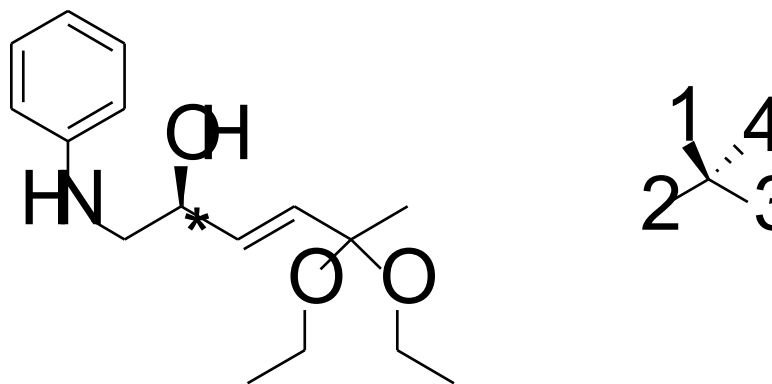
Involved in metabolism

*cis (Z)*

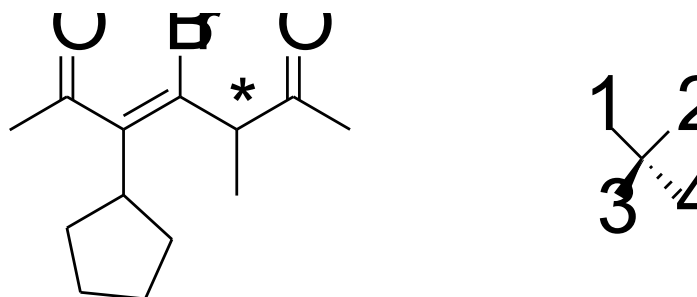


Nakreslete vzorce následujících sloučenin včetně znázornění prostorového uspořádání

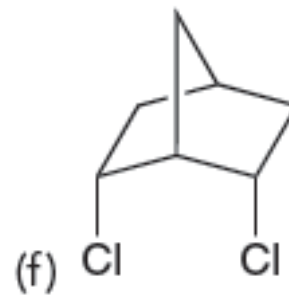
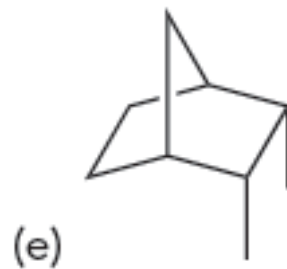
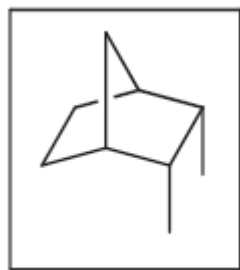
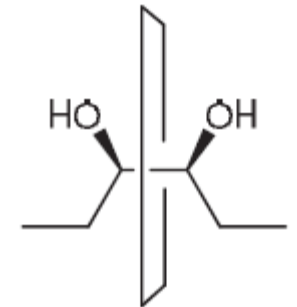
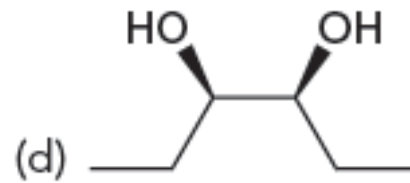
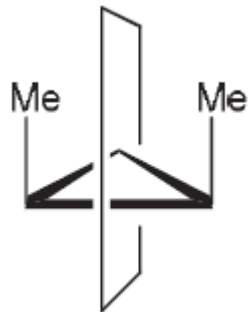
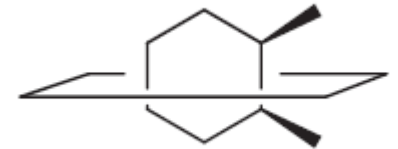
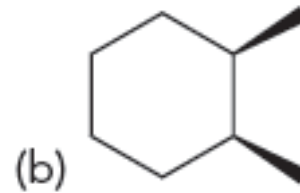
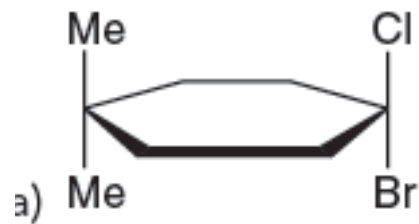
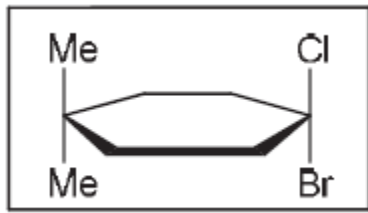
(2*S*,3*E*)-1-(*N*-fenylamino)-5,5-diethoxyhex-3-en-2-ol



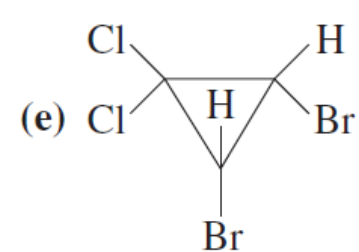
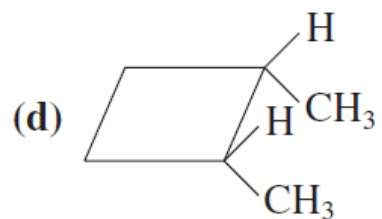
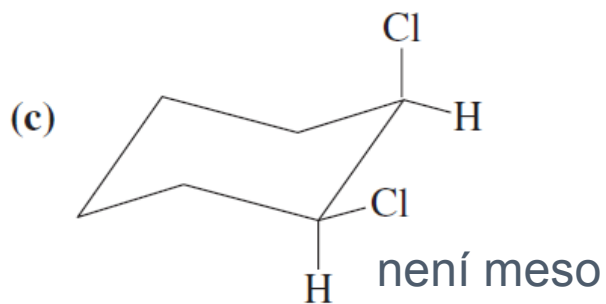
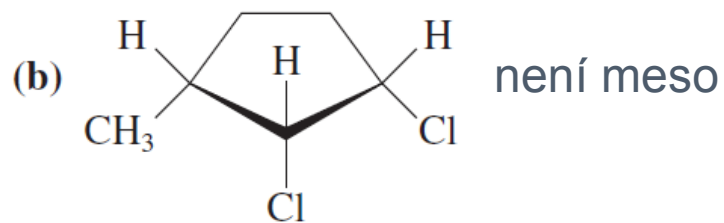
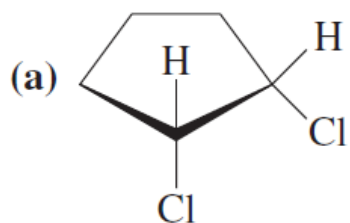
(5*R*,3*Z*)-4-brom-3-cyklopentyl-5-methylhept-3-en-2,6-dion



# Najděte rovinu symetrie u následujících sloučenin

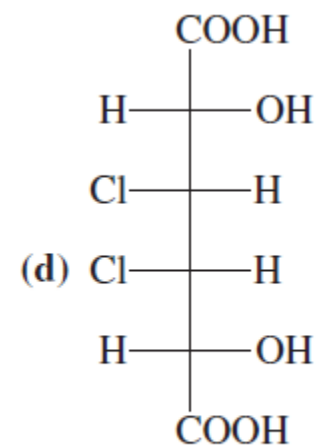
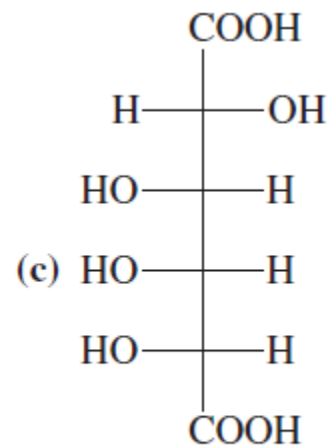
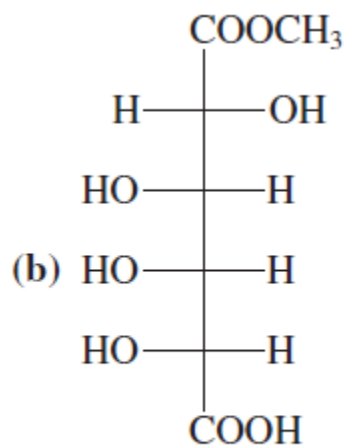
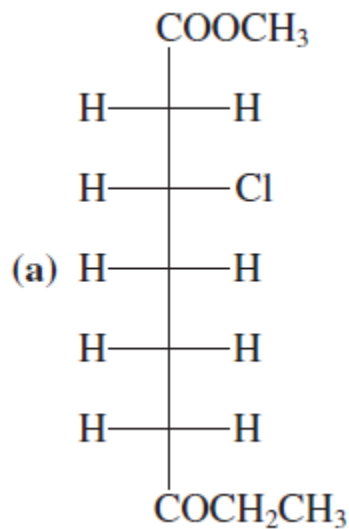


## Která ze sloučenin není *meso* - sloučeninou





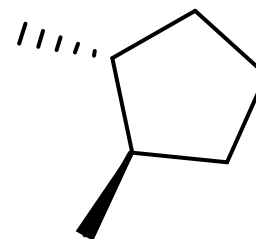
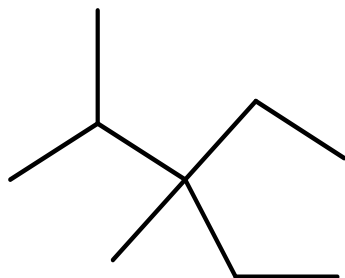
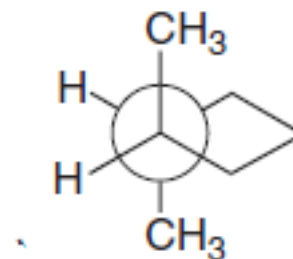
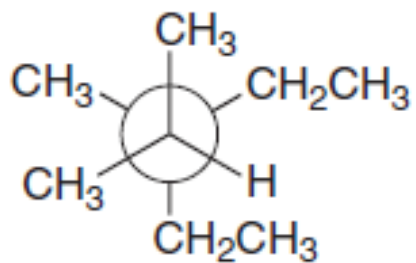
## Která ze sloučenin nebude opticky aktivní



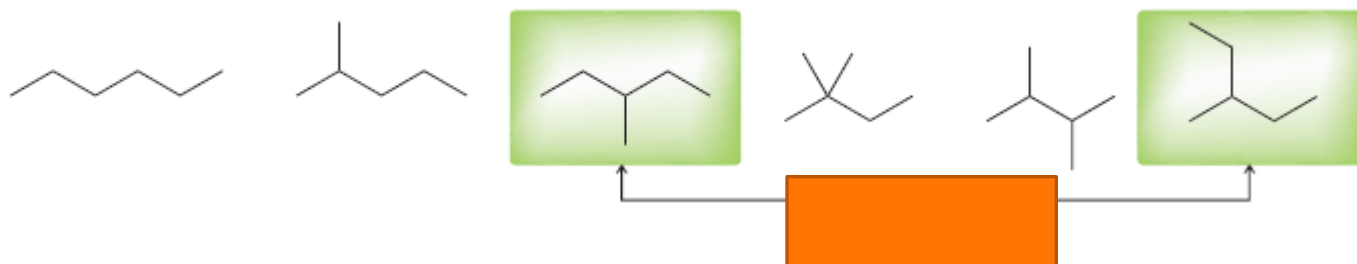
Opticky aktivní není **d**



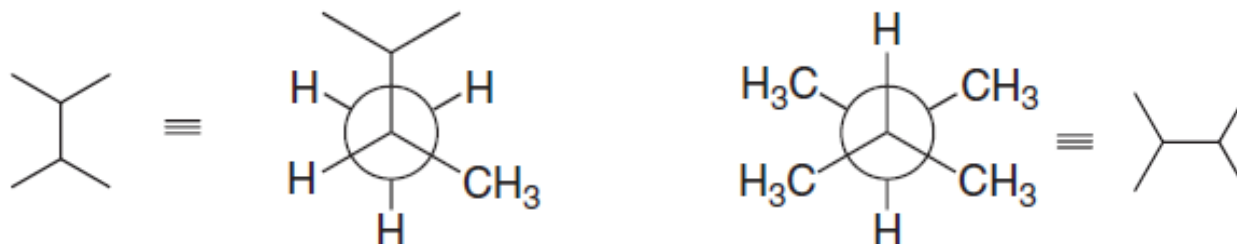
## Nakreslete strukturální vzorce uvedených sloučenin



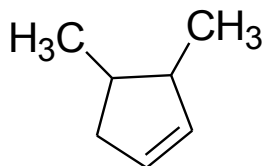
## » Konstituční izomery



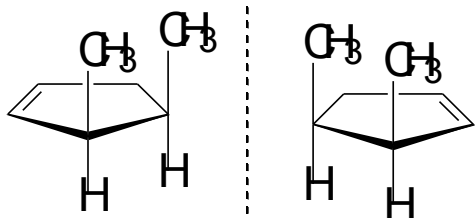
Určete, zda jsou uvedené sloučeniny konstituční izomery



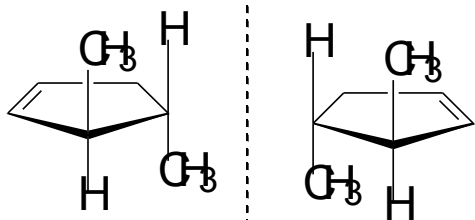
Nakreslete možné stereoizomery uvedených sloučenin. U každé z nich označte, zda otáčí rovinu polarizovaného světla či nikoli.



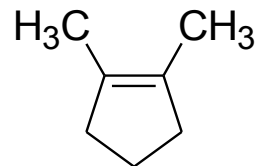
otáčí všechny



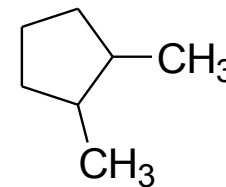
diastereomery



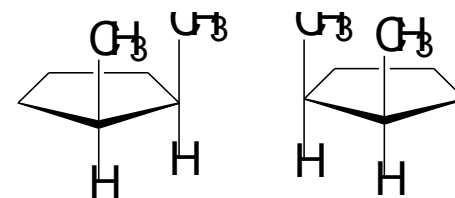
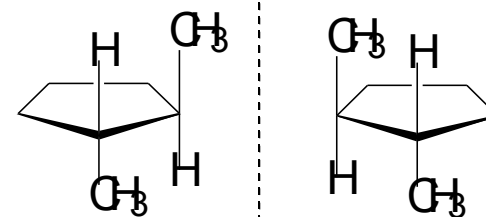
enantiomery



neotáčí



otáčí



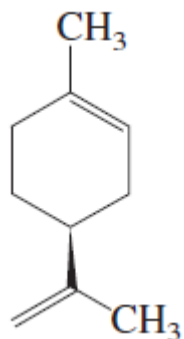
tatožláka  
meso-sloučenina

neotáčí



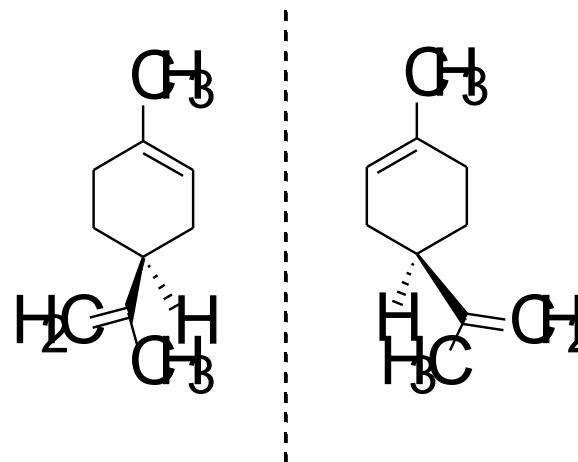
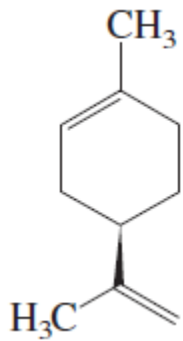
Určete vzájemný stereochemický vztah mezi sloučeninami a pojmenujte je systematickými názvy

(*S*)-limonen

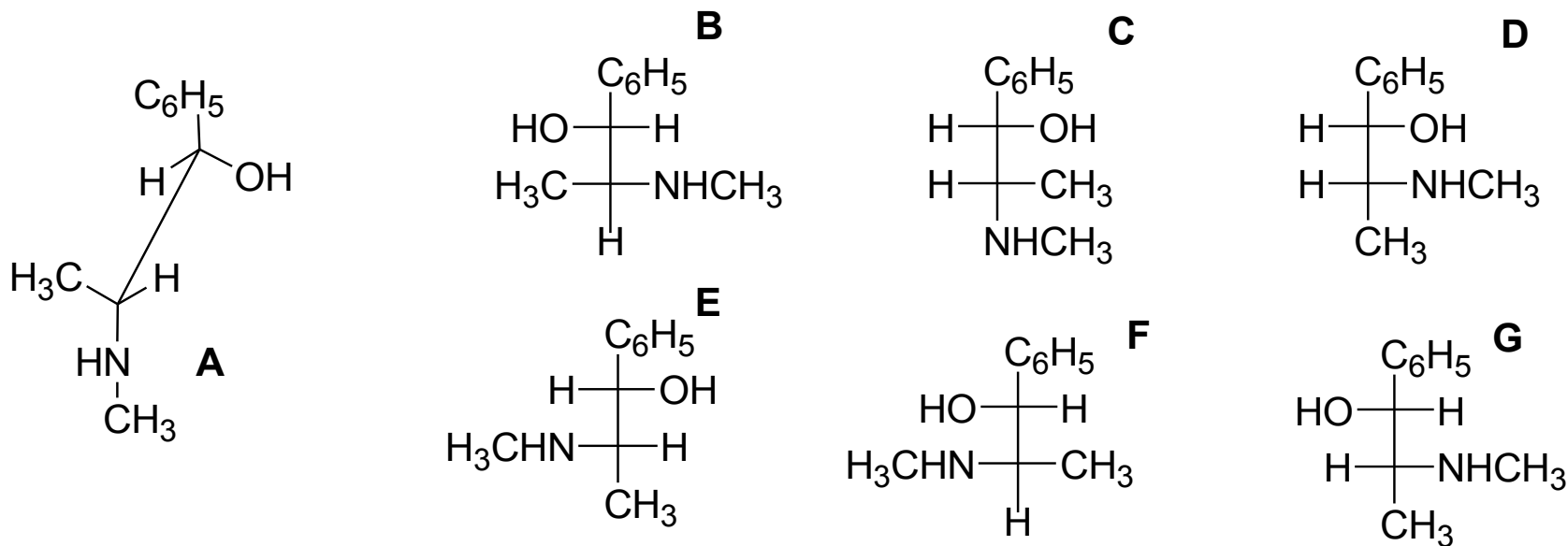


(*S*)-1-methyl-4-(1-methylethenyl)cyclohexen

(*R*)-limonen



Efedrin (A) se používá jako lék, která z uvedených projekcí odpovídá Fischerově projekci látky (A). Vytvořte systematický název efedrinu včetně určení absolutní konfigurace na centrech chiralidy.

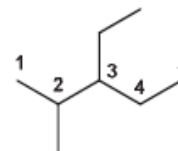
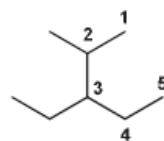
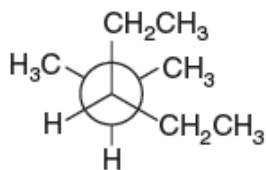
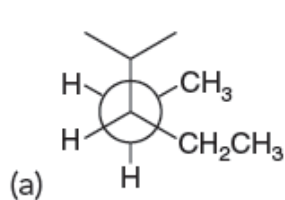


Vzorce **F** a **G**

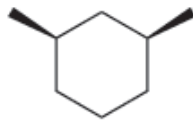
**(1*R*,2*R*)-1-fenyl-2-(*N*-methylamino)propanol**



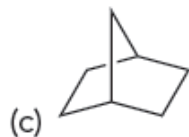
# Určete, jaký je mezi uvedenými sloučeninami vzájemný stereochemický vztah



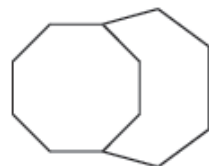
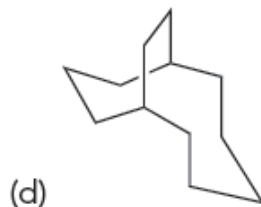
stejné



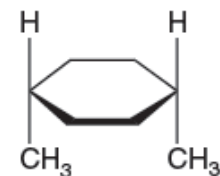
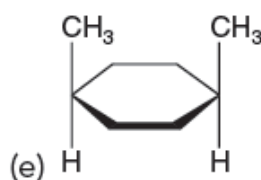
konstituční izomery



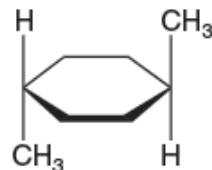
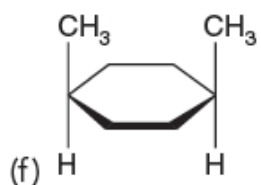
stejné



konstituční izomery

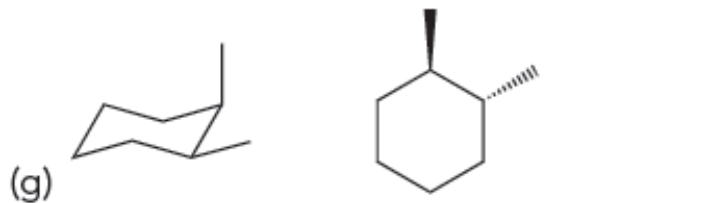


stejné

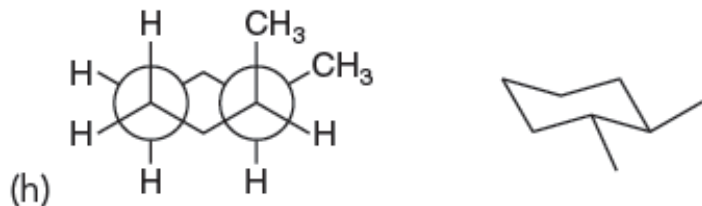


stereoizomery

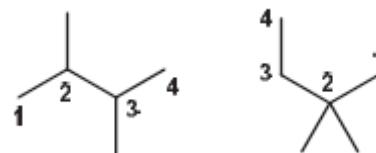
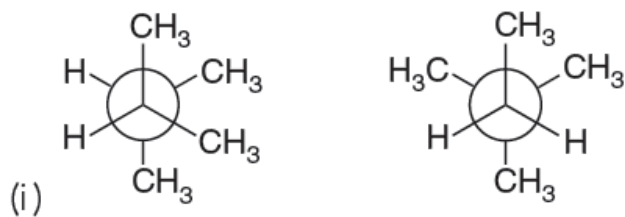




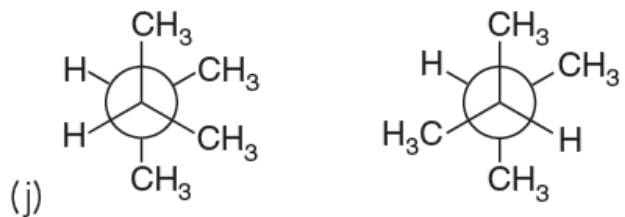
stereoizomery (*cis – trans*, diastereomery)



stereoizomery (*cis – trans*, diastereomery)



konstituční izomery



konformace téže sloučeniny



stereoizomery (*cis – trans*, diastereomery)



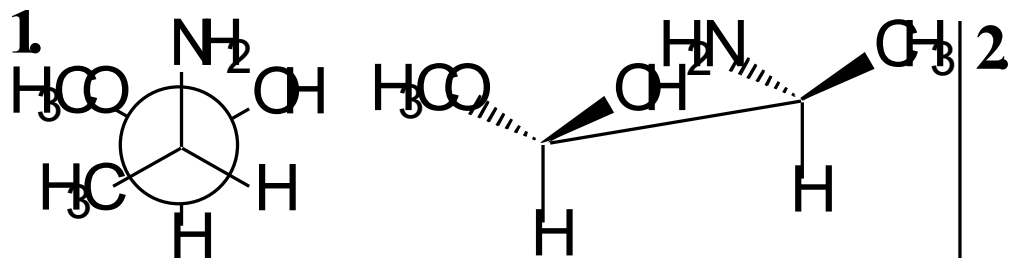
konstituční izomery



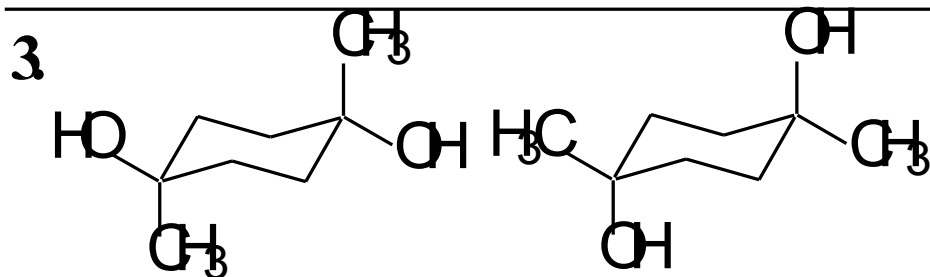
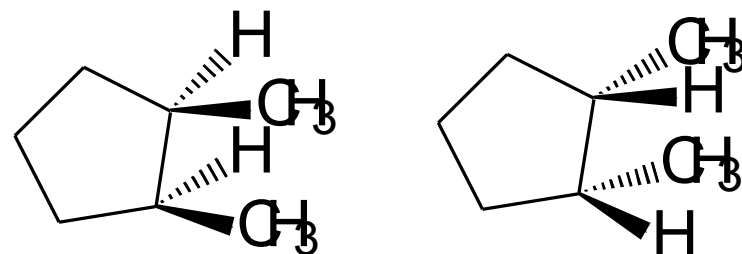


Určete, jaký je mezi uvedenými sloučeninami vzájemný stereochemický vztah

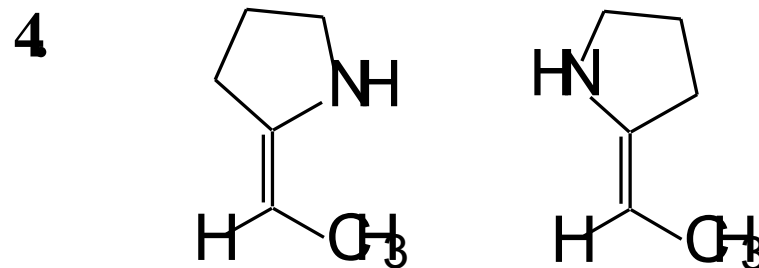
1. diastereomery



2. stejné (*meso*-)



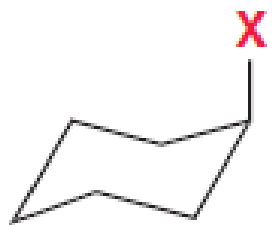
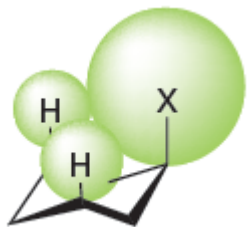
3. konformace téže látky



4. diastereomery  
geometrické izomery



## » Nejstabilnější konformace



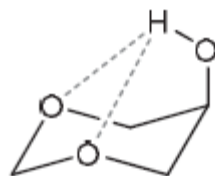
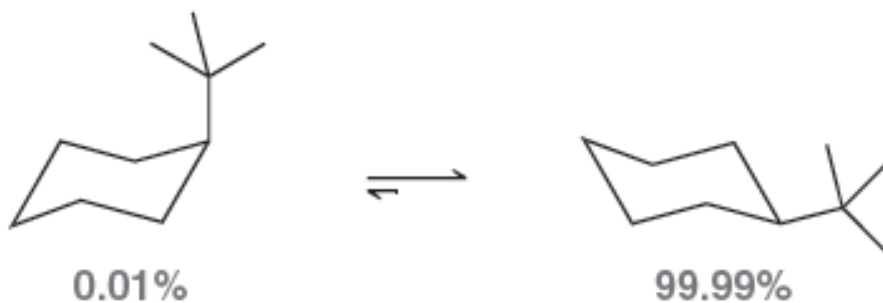
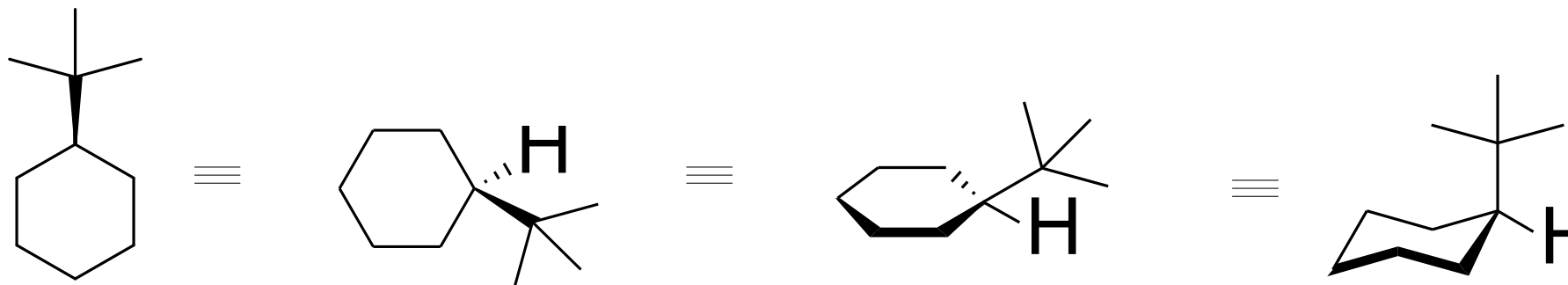
5%



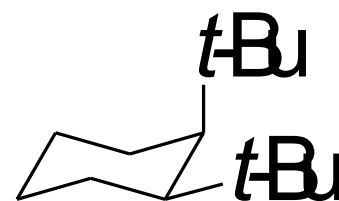
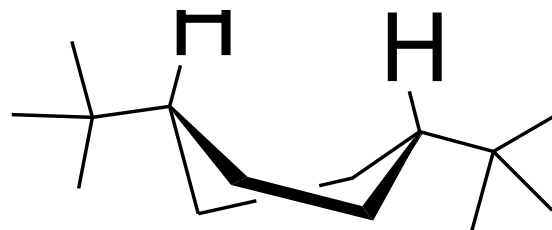
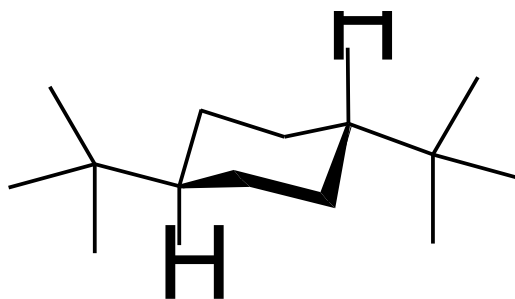
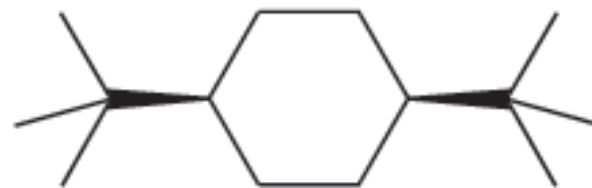
95%



## » Nejstabilnější konformace



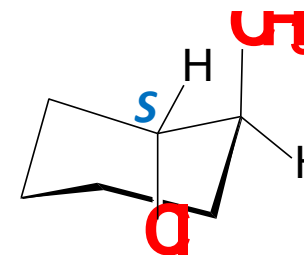
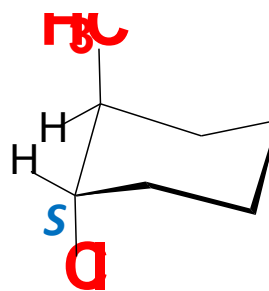
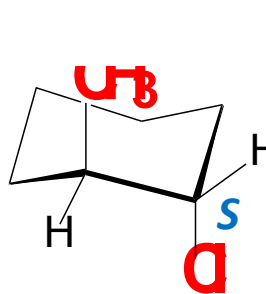
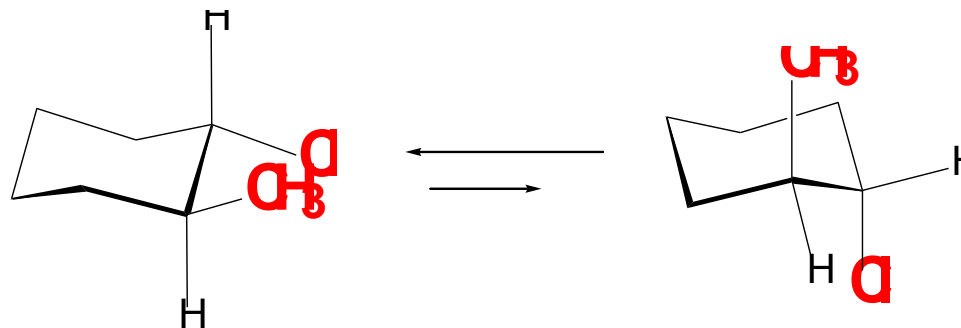
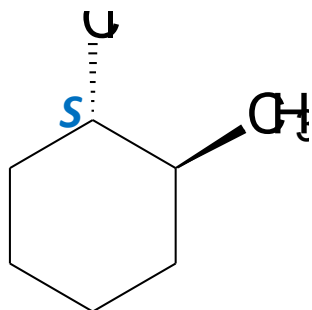
Která ze sloučenin se nachází přednostně v židličkové konformaci a která v twist formě?

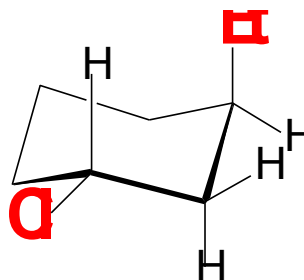
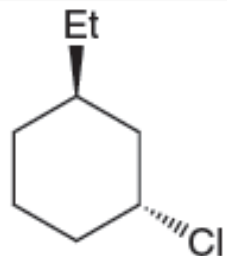
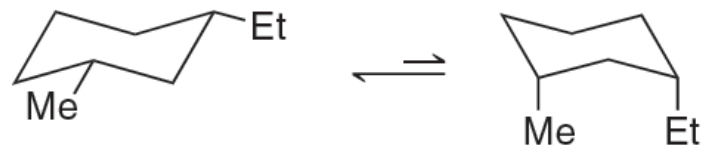
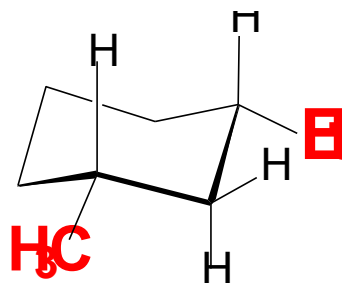
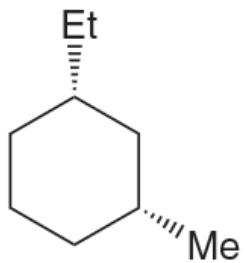


taktore



Uvedenou sloučeninu nakreslete v její nejstabilnější konformaci





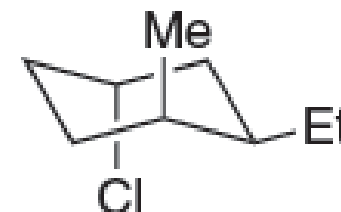
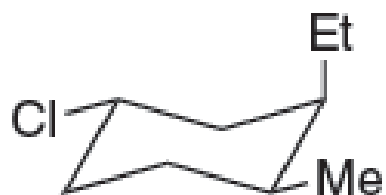
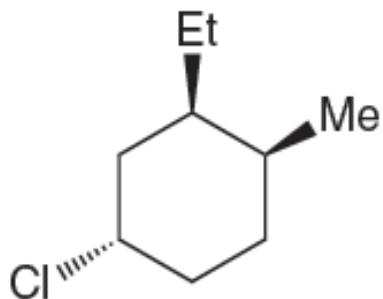
Axial ethyl=8 kJ/mol

Axial chlorine=2 kJ/mol

**TABLE 4.8 1,3-DIAXIAL INTERACTIONS FOR SEVERAL COMMON SUBSTITUENTS**

SUBSTITUENT	1,3-DIAXIAL INTERACTIONS (KJ/MOL)	EQUATORIAL-AXIAL RATIO (AT EQUILIBRIUM)
—Cl	2.0	70 : 30
—OH	4.2	83 : 17
—CH <sub>3</sub>	7.6	95 : 5
—CH <sub>2</sub> CH <sub>3</sub>	8.0	96 : 4
—CH(CH <sub>3</sub> ) <sub>2</sub>	9.2	97 : 3
—C(CH <sub>3</sub> ) <sub>3</sub>	22.8	9999 : 1



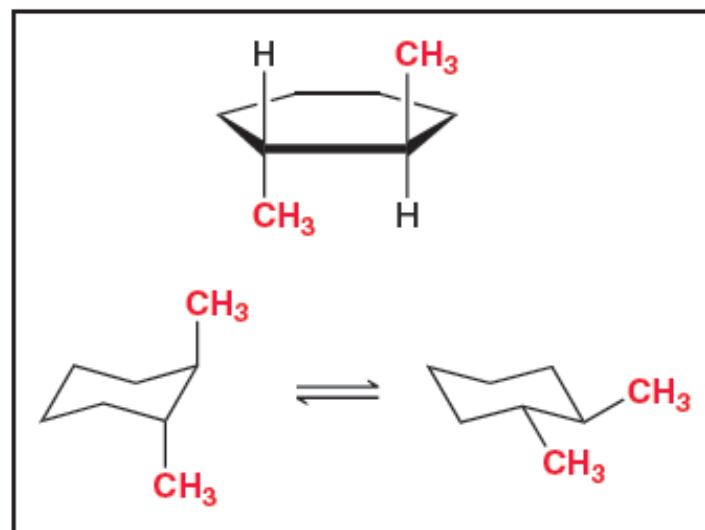
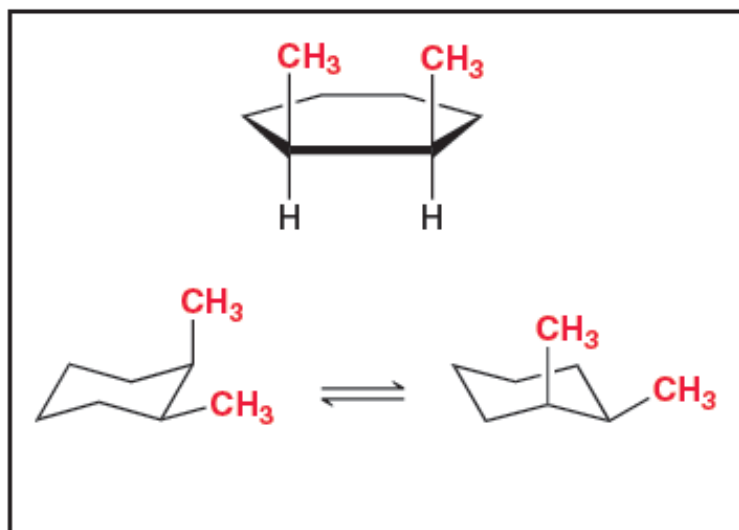


**TABLE 4.8** 1,3-DIAXIAL INTERACTIONS FOR SEVERAL COMMON SUBSTITUENTS

SUBSTITUENT	1,3-DIAXIAL INTERACTIONS (KJ/MOL)	EQUATORIAL-AXIAL RATIO (AT EQUILIBRIUM)
—Cl	2.0	70 : 30
—OH	4.2	83 : 17
—CH <sub>3</sub>	7.6	95 : 5
—CH <sub>2</sub> CH <sub>3</sub>	8.0	96 : 4
—CH(CH <sub>3</sub> ) <sub>2</sub>	9.2	97 : 3
—C(CH <sub>3</sub> ) <sub>3</sub>	22.8	9999 : 1



Určete, který ze dvou možných geometrických izomerů 1,2-dimethylcyklohexanu je termodynamicky stabilnější?



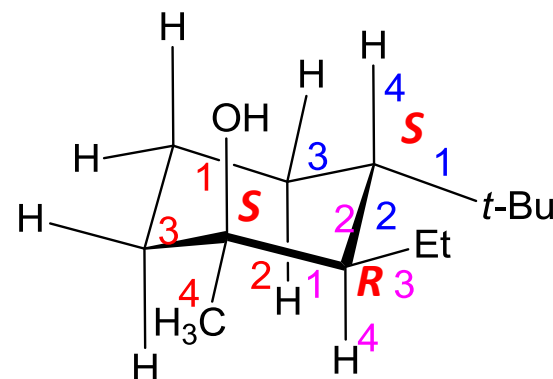
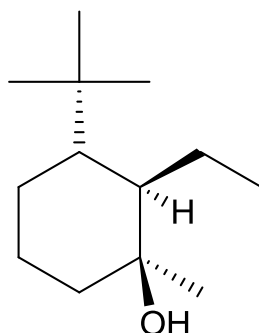
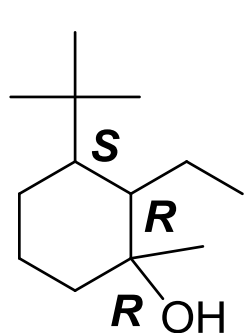
Může zaujmout konformaci, v níž je diekvatoriální pozice obou objemných substituentů



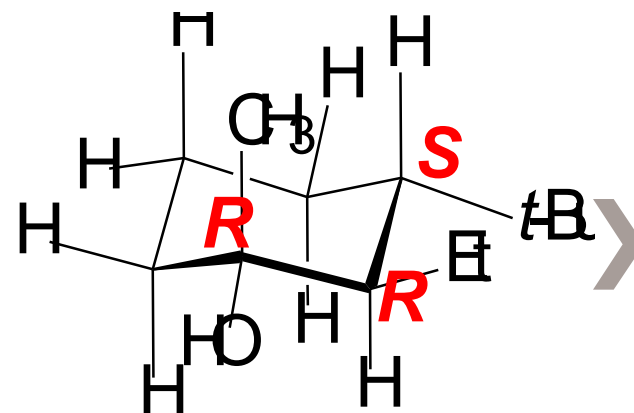


Níže uvedenou sloučeninu nakreslete v její nejstabilnější konformaci

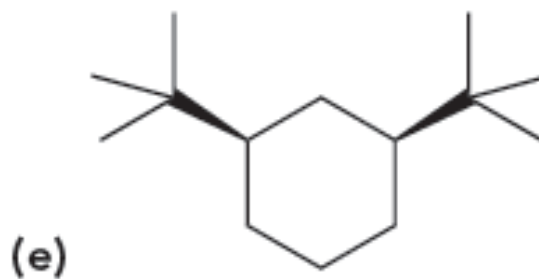
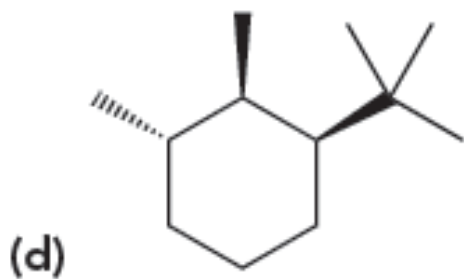
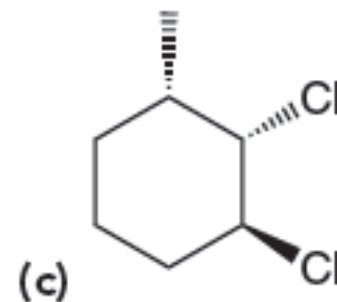
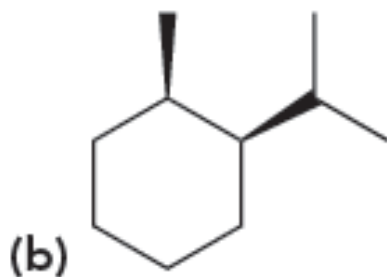
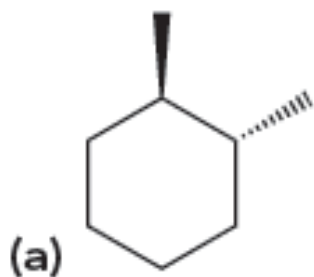
**(1*R*,2*R*,3*S*)-3-*tert*-butyl-2-ethyl-1-methylcyklohexanol**



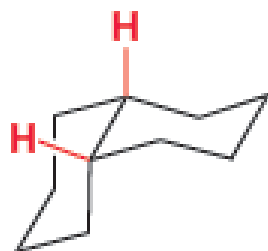
- 1) Nakreslíme strukturní vzorec sloučeniny
- 2) Libovolně umístíme substituenty, *t*-Bu umístíme do ekvatoriální pozice, neboť nejstabilnější konformace preferuje objemný substituent v ekv. pozici
- 3) Určíme abs. konfiguraci
- 4) Na uhlíku C1 prohodíme dva substituenty, neboť dle zadání má být abs. konfigurace *R* a nám vyšla *S*



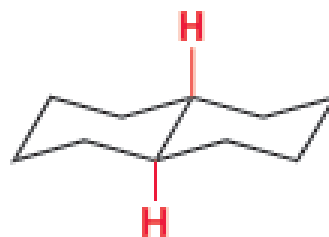
# Nakreslete nejstabilnější konformaci uvedených sloučenin



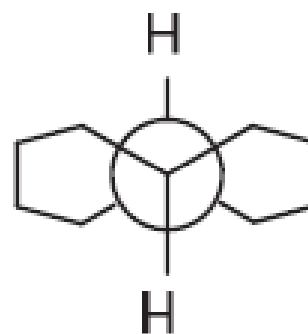
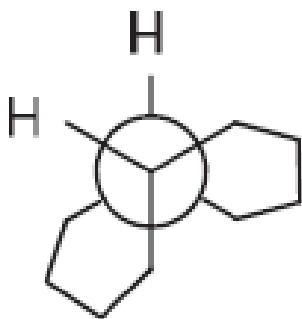
Určete, který ze dvou možných geometrických izomerů dekalinu je termodynamicky stabilnější?



*cis*-Decalin



*trans*-Decalin

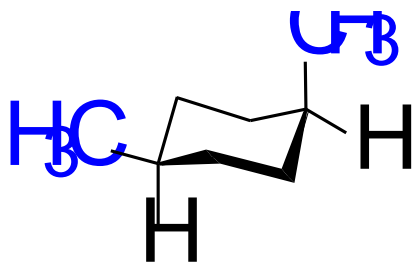


*trans*- dekalin je stabilnější – pouze 2 gauche interakce (*cis* má 3)



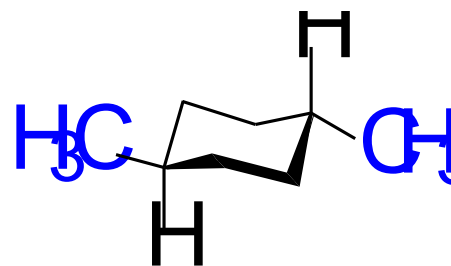
Určete, který z izomerů dimethylcyklohexanu je chirální?

1,4-dimethylcyklohexan



*cis*

achirální, rovina symetrie



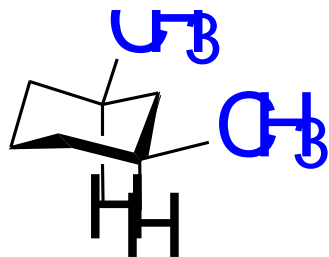
*trans*

achirální, rovina symetrie



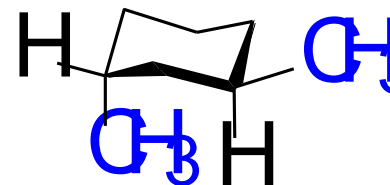
Určete, který z izomerů dimethylcyklohexanu je chirální?

1,3-dimethylcyklohexan



*cis*

achirální, rovina symetrie



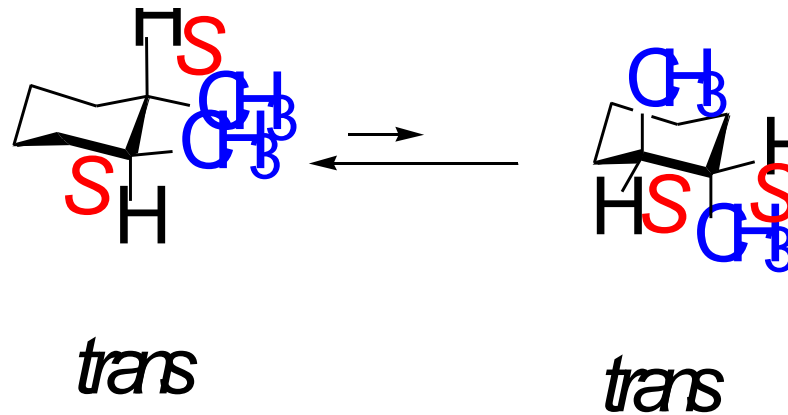
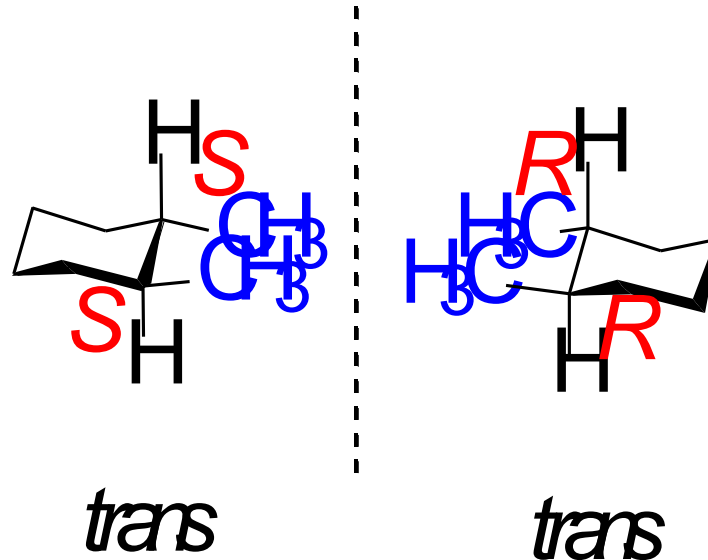
*trans*

chirální



Určete, který z izomerů dimethylcyklohexanu je chirální?

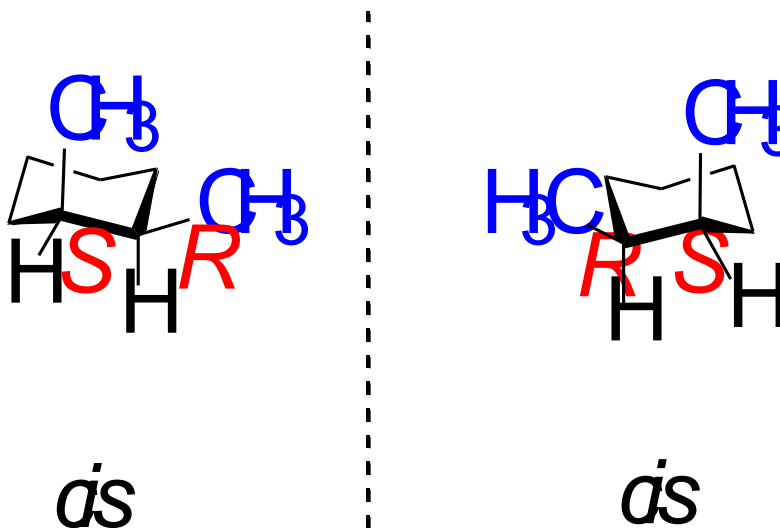
1,2-dimethylcyklohexan



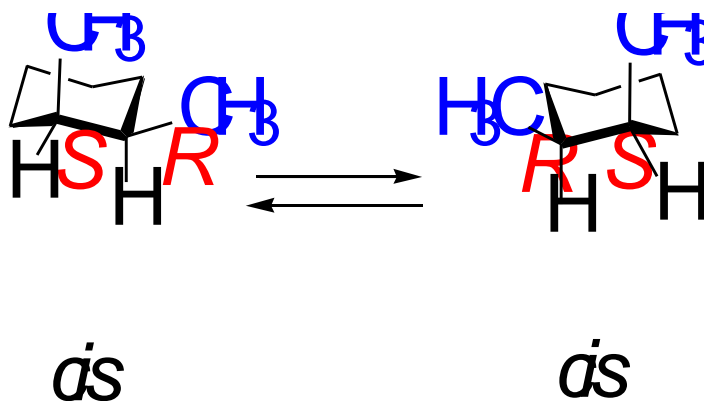
chirální

# Určete, který z izomerů dimethylcyklohexanu je chirální?

1,2-dimethylcyklohexan



chirální **ALE POZOR!**



za normální teploty je není možné separovat a jedná se o konformace téže látky

