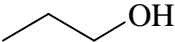

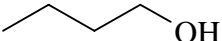
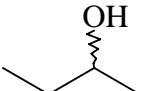
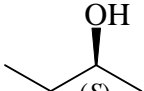
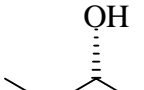
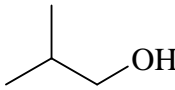
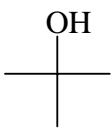


Homologe Reihe der unverzweigten primären Alkohole  $C_nH_{(2n+2)}O$ 

systematischer Name	Konstitution	Siedepunkt [°C]	Schmelzpunkt [°C]
Methanol	$CH_3-OH$	65	- 98
Ethanol	$CH_3-CH_2-OH$	78	- 117
1-Propanol	$CH_3-CH_2-CH_2-OH$	97	- 126
1-Butanol	$CH_3-(CH_2)_2-CH_2-OH$	118	- 89
1-Pentanol	$CH_3-(CH_2)_3-CH_2-OH$	138	- 78
1-Hexanol	$CH_3-(CH_2)_4-CH_2-OH$	157	- 52
1-Heptanol	$CH_3-(CH_2)_5-CH_2-OH$	175	- 36
1-Octanol	$CH_3-(CH_2)_6-CH_2-OH$	194	- 15
⋮	⋮	⋮	⋮
1-Hexandecanol	$CH_3-(CH_2)_{14}-CH_2-OH$	344	50

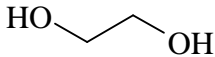
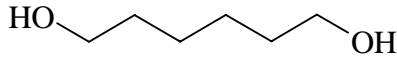
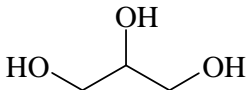
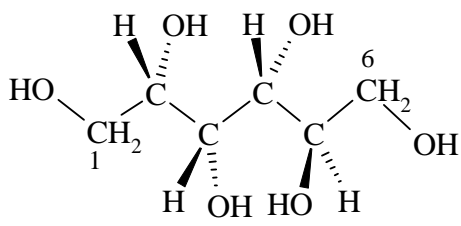
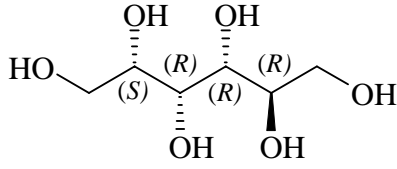
## Konstitutions- und konfigurationsisomere Alkohole

systematischer Name (Trivialname)	ausführliche Schreibweise	Stenographie	Siedep. [°C]	Schmelzp. [°C]
<b>1-Propanol</b> (n-Propylalkohol)	$\text{CH}_3\text{-CH}_2\text{-CH}_2\text{-OH}$		97	-126
<b>2-Propanol</b> (Isopropylalkohol)	$\begin{array}{c} \text{OH} \\   \\ \text{CH}_3\text{-CH-CH}_3 \end{array}$		82	-89
<b>1-Butanol</b> (n-Butylalkohol)	$\text{CH}_3\text{-CH}_2\text{-CH}_2\text{-CH}_2\text{-OH}$		118	-89
<b>2-Butanol</b> (sekundärer Butylalkohol) ( <i>sek</i> -Butanol)	$\begin{array}{c} \text{OH} \\   \\ \text{CH}_3\text{-CH}_2\text{-CH-CH}_3 \end{array}$		100	-114
	$\begin{array}{c} \text{HO} \quad \text{H} \\ \diagdown \quad \diagup \\ \text{H}_3\text{C-CH}_2\text{-C-CH}_3 \end{array}$			
	$\begin{array}{c} \text{H} \quad \text{OH} \\ \diagdown \quad \diagup \\ \text{H}_3\text{C-CH}_2\text{-C-CH}_3 \end{array}$			
<b>2-Methyl-1-propanol</b> (Isobutylalkohol)	$\begin{array}{c} \text{CH}_3 \\   \\ \text{H}_3\text{C-CH-CH}_2\text{-OH} \end{array}$		108	-108
<b>2-Methyl-2-propanol</b> (tertiärer Butylalkohol) ( <i>tert</i> -Butanol)	$\begin{array}{c} \text{OH} \\   \\ \text{CH}_3\text{-C-CH}_3 \\   \\ \text{CH}_3 \end{array}$		82	24

## Vergleich unverzweigter Verbindungen mit ähnlicher Molmasse

systematischer Name ausführliche Schreibweise	Stenographie	Mol- masse	Dipol- moment [D]	Siede- punkt [°C]	Schmelz- punkt [°C]
Pentan $\text{CH}_3\text{—CH}_2\text{—CH}_2\text{—CH}_2\text{—CH}_3$		72	0	36	– 129
Diethylether $\text{CH}_3\text{—CH}_2\text{—O—CH}_2\text{—CH}_3$		74	1.18	35	– 116
1-Chlorpropan $\text{CH}_3\text{—CH}_2\text{—CH}_2\text{—Cl}$		79	2.10	47	– 123
1-Butanal $\text{CH}_3\text{—CH}_2\text{—CH}_2\text{—C}\begin{matrix} \text{O} \\ \parallel \\ \text{H} \end{matrix}$		72	2.72	75	– 96
1-Butanol $\text{CH}_3\text{—CH}_2\text{—CH}_2\text{—CH}_2\text{—OH}$		74	1.63	118	– 90
Propansäure $\text{CH}_2\text{—CH}_2\text{—C}\begin{matrix} \text{O} \\ \parallel \\ \text{OH} \end{matrix}$		74	1.68	141	– 22

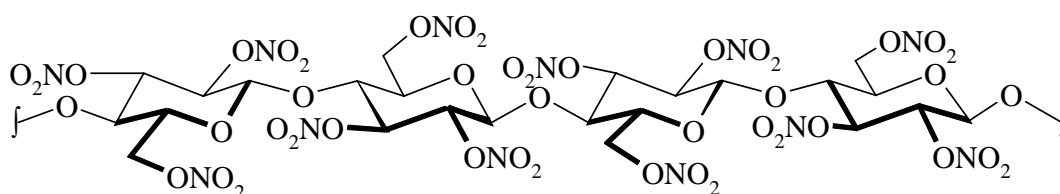
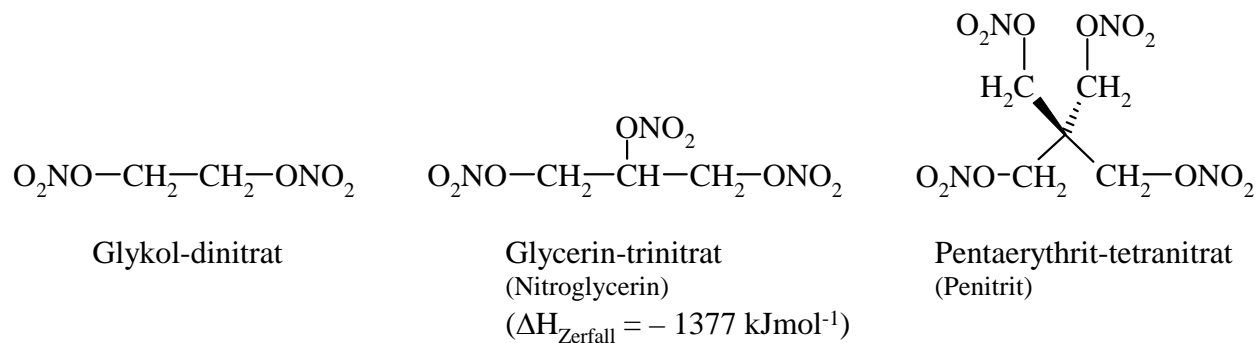
## Alkohole mit mehreren Hydroxygruppen

systematischer Name (Trivialname)	Struktur		Siedepunkt [°C]	Schmelzpunkt [°C]
	ausführliche Schreibweise	Stenographie		
1,2-Ethandiol (Ethylenglykol) (Glykol)	HO-CH <sub>2</sub> -CH <sub>2</sub> -OH		197	-12
1,6-Hexandiol	HO-CH <sub>2</sub> -CH <sub>2</sub> -CH <sub>2</sub> -CH <sub>2</sub> -CH <sub>2</sub> -CH <sub>2</sub> -OH		250	42
1,2,3-Propantriol (Glycerin)	$\begin{array}{c} \text{OH} \\   \\ \text{HO}-\text{CH}_2-\text{CH}-\text{CH}_2-\text{OH} \end{array}$		290	18
1,2,3,4,5,6-Hexanhexol (D-Sorbit)				92

## Mineralsäureester

Mineralsäure		Alkylester der Mineralsäure (R = Alkyl)	
Salpetersäure $\text{HNO}_3$	$\text{O}_2\text{N}-\text{OH}$	$\text{O}_2\text{N}-\text{O}-\text{R}$	Salpetersäurealkylester (Alkylnitrate)
Schwefelsäure $\text{H}_2\text{SO}_4$	$\begin{array}{c} \text{OH} \\ \diagup \\ \text{O}_2\text{S} \\ \diagdown \\ \text{OH} \end{array}$	$\begin{array}{c} \text{O}-\text{R} \\ \diagup \\ \text{O}_2\text{S} \\ \diagdown \\ \text{OH} \end{array}$	Schwefelsäuremonoalkylester (Alkylsulfate)
		$\begin{array}{c} \text{O}-\text{R} \\ \diagup \\ \text{O}_2\text{S} \\ \diagdown \\ \text{O}-\text{R} \end{array}$	Schwefelsäuredialkylester (Disulfate)
Phosphorsäure $\text{H}_3\text{PO}_4$	$\begin{array}{c} \text{OH} \\ \diagup \\ \text{OP} \\ \diagdown \\ \text{OH} \\ \text{OH} \end{array}$	$\begin{array}{c} \text{O}-\text{R} \\ \diagup \\ \text{OP} \\ \diagdown \\ \text{OH} \\ \text{OH} \end{array}$	Phosphorsäuremonoalkylester (Alkylphosphate)
		$\begin{array}{c} \text{O}-\text{R} \\ \diagup \\ \text{OP} \\ \diagdown \\ \text{O}-\text{R} \\ \text{OH} \end{array}$	Phosphorsäuredialkylester (Dialkylphosphate)
		$\begin{array}{c} \text{O}-\text{R} \\ \diagup \\ \text{OP} \\ \diagdown \\ \text{O}-\text{R} \\ \text{O}-\text{R} \end{array}$	Phosphorsäuretrialkylester (Trialkylphosphate)

### Salpetersäureester von Polyalkoholen sind Explosivstoffe



Cellulose-trinitrat  
(Schießbaumwolle)

## Alkoholgehalt von Getränken

Getränk	Frucht	Stammwürze [%]	Ethanol [Vol%]
Bier	Gerste	11 – 14	4.6 – 5.8
Starkbier		> 16	> 6.7
Apfelwein	Äpfel		8 – 9
Wein	Trauben		8 – 13
Liköre	Diverse Früchte		20 – 60
Branntwein	Diverse Früchte		32 – 60


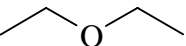
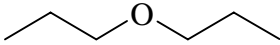
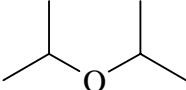
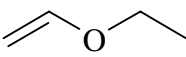
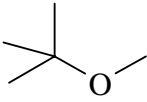
Quelle: AID-Verbraucherdienst 35 (1990) Heft 5, S. 9

## Flüchtige Alkohole außer Ethanol im Wein

Alkohole	Weißwein (mg/l)	Rotwein (mg/l)
Methanol	38 – 118	43 – 222
1-Propanol	9 – 48	11 – 52
2-Methyl-1-propanol	28 – 170	45 – 140
1-Butanol	1.4 – 8.5	2.1 – 2.3
(S)-(-)-2-Methyl-1-butanol	17 – 82	48 – 150
3-Methyl-1-butanol	70 – 320	117 – 490
1-Hexanol	3 – 10	3 – 10

Quelle: H.-D. Belitz, W. Grosch, Lehrbuch der Lebensmittelchemie, 3. Aufl., Springer, Heidelberg-Berlin, 1987

## Acyclische Ether

systematischer Name (Trivialname)	Konstitution		Siedep. [°C]
	ausführliche Schreibweise	Stenographie	
Dimethylether	$\text{H}_3\text{C}-\text{O}-\text{CH}_3$		- 23
Diethylether (Ether)	$\text{H}_3\text{C}-\text{CH}_2-\text{O}-\text{CH}_2-\text{CH}_3$		35
Di-1-propylether	$\text{H}_3\text{C}-\text{CH}_2-\text{CH}_2-\text{O}-\text{CH}_2-\text{CH}_2-\text{CH}_3$		91
Di-2-propylether	$\begin{array}{c} \text{CH}_3 \quad \text{CH}_3 \\   \quad   \\ \text{H}_3\text{C}-\text{CH}-\text{O}-\text{CH}-\text{CH}_3 \end{array}$		68
Ethylvinylether	$\text{H}_2\text{C}=\text{CH}-\text{O}-\text{CH}_2-\text{CH}_3$		36
tert-Butylmethyl- ether (MtB)	$\begin{array}{c} \text{CH}_3 \\   \\ \text{H}_3\text{C}-\text{C}-\text{O}-\text{CH}_3 \\   \\ \text{CH}_3 \end{array}$		55



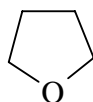
### Cyclische Ether (das Sauerstoffatom ist Teil eines Ringes)



**Oxiran**  
(Ethylenoxid)

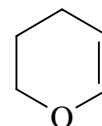
Siedepunkt [°C]

14



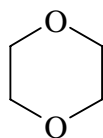
**Tetrahydrofuran**  
(THF)

67



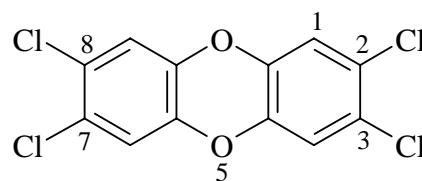
**2,3-Dihydropyran**

86



**1,4-Dioxan**

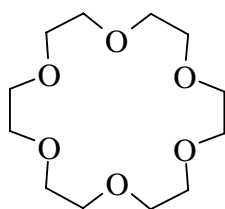
MAK = 180 mg/m<sup>3</sup>



**2,3,7,8-Tetrachlordibenzo-*p*-dioxin**  
(TCDD)

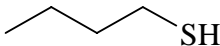
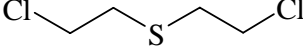
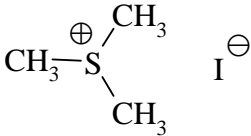
LD<sub>50</sub> = 0.001 mg/kg Meerschweinchen

LD<sub>50</sub> = 3 mg/kg Hamster


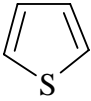

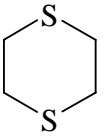


**18-Krone-6**  
(ein Kronenether)

## Acyclische Schwefel(II)-Verbindungen

Name (Trivialname)	Konstitution		Siedep. [°C]
	ausführliche Schreibweise	Stenographie	
<b>Methanthiol</b> (Methylmercaptan)	$\text{H}_3\text{C}-\text{SH}$		6
<b>1-Butanthiol</b> ( <i>n</i> -Butylmercaptan)	$\text{H}_3\text{C}-\text{CH}_2-\text{CH}_2-\text{CH}_2-\text{SH}$		98
<b>Dimethylthioether</b> (Dimethylsulfid)	$\text{H}_3\text{C}-\text{S}-\text{CH}_3$		37
<b>S-Lost</b> (chemischer Kampfstoff, cancerogen)	$\text{Cl}-\text{CH}_2-\text{CH}_2-\text{S}-\text{CH}_2-\text{CH}_2-\text{Cl}$		217
<b>Trimethylsulfoniumiodid</b>			
<b>Dimethyldisulfid</b>	$\text{CH}_3-\text{S}-\text{S}-\text{CH}_3$		110

## Cyclische Schwefel(II)-Verbindungen (das Schwefelatom ist Teil des Ringes)

			
Thiiran	Thiophen	1,3-Dithian	1,4-Dithian
Siedepunkt [°C] 55	84	207	199

## 8.11

## Schwefel(II)-haltige Carbonsäuren

