

Cyclohexanol, 5-methyl-2-(1-methylethyl)-, (1 «alpha»,2 «beta»,5 «alpha»)-

Other names:

Menthol, cis-1,3,trans-1,4-

Hexahydrothymol

Menthacamphor

Menthol

Menthomenthol

Peppermint camphor

dl-Menthol

p-Menthan-3-ol

Cyclohexanol, 2-isopropyl-5-methyl-

NCI-C50000

5-Methyl-2-(1-methylethyl)cyclohexanol

2-Isopropyl-5-methylcyclohexanol

l-Menthol

3-p-Menthanol

5 «alpha»-Methyl-2 «beta»-(1 «alpha»-methylethyl)cyclohexanol

component of Dermoplast

Fancol menthol

Fisherman's friend lozenges

Menthol terpine hydrate

component of Minut-rub

Racemic menthol

component of Robitussin cough drops

component of Sarna

component of Theragesic

Therapeutic mineral ice

Cyclohexanol, 2-isopropyl-5-methyl-, (1 «alpha»,2 «beta»,5 «alpha»)-

2-Isopropyl-5-methylcyclohexanol, (1 «alpha»,2 «beta»,5 «alpha»)-

5-Methyl-2-(1-methylethyl)cyclohexanol, (1 «alpha»,2 «beta»,5 «alpha»)-

(. +/-)-Menthol

Cyclohexanol, 5-methyl-2-(1-methylethyl)-, (1R,2S,5R)-rel-

NSC 2603

Racementhol

rac-Menthol

Diterpene

Cyclohexanol, 5-methyl-2-(1-methylethyl)-, (1 «alpha»,2 «beta»,5 «alpha»)-(. +/-)-

Inchi: InChI=1S/C10H20O/c1-7(2)9-5-4-8(3)6-10(9)11/h7-11H,4-6H2,1-3H3/t8-,9+,10-/m1/s1

InchiKey: NOOLISFMXDJSKH-KXUCPTDWSA-N

Formula: C10H20O

SMILES: CC1CCC(C(C)C)C(O)C1

Mol. weight [g/mol]: 156.27

Physical Properties

Property code	Value	Unit	Source
gf	-96.91	kJ/mol	Joback Method
hf	-393.60	kJ/mol	Joback Method
h _{fus}	16.20	kJ/mol	Joback Method
h _{vap}	53.96	kJ/mol	Joback Method
log ₁₀ ws	-2.55		Crippen Method
logp	2.440		Crippen Method
m _{cvol}	146.770	ml/mol	McGowan Method
pc	2659.77	kPa	Joback Method
tb	530.15	K	Joback Method
tc	719.67	K	Joback Method
tf	247.18	K	Joback Method
vc	0.539	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	370.36	J/mol×K	530.15	Joback Method
cpg	387.84	J/mol×K	561.74	Joback Method
cpg	404.52	J/mol×K	593.32	Joback Method
cpg	420.40	J/mol×K	624.91	Joback Method
cpg	435.50	J/mol×K	656.49	Joback Method
cpg	449.82	J/mol×K	688.08	Joback Method
cpg	463.40	J/mol×K	719.67	Joback Method
d _{visc}	0.0430815	Paxs	247.18	Joback Method
d _{visc}	0.0077902	Paxs	294.34	Joback Method
d _{visc}	0.0022592	Paxs	341.50	Joback Method
d _{visc}	0.0008847	Paxs	388.66	Joback Method
d _{visc}	0.0004244	Paxs	435.83	Joback Method
d _{visc}	0.0002350	Paxs	482.99	Joback Method
d _{visc}	0.0001446	Paxs	530.15	Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	376.70	K	1.00	NIST Webbook

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C15356704&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tbrp:	Boiling point at reduced pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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