(1S,2R,5R)-2-Methyl-5-((R)-6-methylhept-5-en-2-yl)

Other names: Bicyclo[3.1.0]hexan-2-ol, 5-[(1R)-1,5-dimethyl-4-hexen-1-yl]-2-methyl-, (1S,2R,5R)-
Bicyclo[3.1.0]hexan-2-ol, 5-[(1R)-1,5-dimethyl-4-hexenyl]-2-methyl-, (1S,2R,5R)-
cis-Sesquisabinene hydrate
(Z)-«beta»-Sesquisabinene hydrate
(Z)-sesquisabinenehydrate
cis-Sesquisabinenhydrate
Sesquisabinene hydrate (cis)

Inchi: InChI=1S/C15H26O/c1-11(2)6-5-7-12(3)15-9-8-14(4,16)13(15)10-15/h6,12-13,16H,5,7-1

InchiKey: IRDFGGRWKUKANK-KBUPBQIOSA-N

Formula: C15H26O

SMILES: CC(C)=CCCC(C)C12CCC(C)(O)C1C2

Mol. weight [g/mol]: 222.37

CAS: 58319-05-4

Physical Properties

Property code	Value	Unit	Source
gf	110.64	kJ/mol	Joback Method
hf	-247.27	kJ/mol	Joback Method
hfus	18.81	kJ/mol	Joback Method
hvap	62.53	kJ/mol	Joback Method
log10ws	-4.39		Crippen Method
logp	3.920		Crippen Method
mcvol	202.060	ml/mol	McGowan Method
pc	2096.50	kPa	Joback Method
rinpol	1542.00		NIST Webbook
rinpol	1543.00		NIST Webbook
rinpol	1545.00		NIST Webbook
rinpol	1590.00		NIST Webbook
rinpol	1567.00		NIST Webbook
rinpol	1525.00		NIST Webbook
rinpol	1565.00		NIST Webbook
rinpol	1536.00		NIST Webbook
rinpol	1560.00		NIST Webbook
rinpol	1540.00		NIST Webbook
rinpol	1539.00		NIST Webbook
rinpol	1554.00		NIST Webbook
rinpol	1543.00		NIST Webbook

rinpol	1521.00		NIST Webbook
rinpol	1524.00		NIST Webbook
rinpol	1544.00		NIST Webbook
rinpol	1562.00		NIST Webbook
rinpol	1542.00		NIST Webbook
rinpol	1544.00		NIST Webbook
rinpol	1545.00		NIST Webbook
rinpol	1581.00		NIST Webbook
rinpol	1572.00		NIST Webbook
rinpol	1539.00		NIST Webbook
rinpol	1540.00		NIST Webbook
rinpol	1530.00		NIST Webbook
rinpol	1540.00		NIST Webbook
ripol	2075.00		NIST Webbook
ripol	2088.00		NIST Webbook
tb	647.67	K	Joback Method
tc	844.14	K	Joback Method
tf	365.03	K	Joback Method
vc	0.778	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	583.77	J/mol×K	647.67	Joback Method
cpg	600.91	J/mol×K	680.41	Joback Method
cpg	617.33	J/mol×K	713.16	Joback Method
cpg	633.24	J/mol×K	745.90	Joback Method
cpg	648.84	J/mol×K	778.65	Joback Method
cpg	664.35	J/mol×K	811.39	Joback Method
cpg	679.97	J/mol×K	844.14	Joback Method

Sources

Joback Method: https://en.wikipedia.org/wiki/Joback_method

McGowan Method: http://link.springer.com/article/10.1007/BF02311772

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C58319054&Units=SI

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cpg:	Ideal gas heat capacity
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
rinpol:	Non-polar retention indices
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume