

Isopulegol acetate

Other names:	Cyclohexanol, 5-methyl-2-(1-methylethenyl)-, acetate, [1R-(1«alpha»,2«beta»,5«alpha»)]-isopulegyl acetate Cyclohexanol, 5-methyl-2-(1-methylethenyl)-, 1-acetate, (1R,2S,5R)-(-)-Isopulegyl acetate Isopulegol acetate (Isomer 2) isopulegol acetate (isomer 1)
Inchi:	InChI=1S/C12H20O2/c1-8(2)11-6-5-9(3)7-12(11)14-10(4)13/h9,11-12H,1,5-7H2,2-4H3/t9
InchiKey:	HLHIVJRLODSUCI-ADEWGFFLSA-N
Formula:	C12H20O2
SMILES:	C=C(C)C1CCC(C)CC1OC(C)=O
Mol. weight [g/mol]:	196.29
CAS:	57576-09-7

Physical Properties

Property code	Value	Unit	Source
gf	-95.44	kJ/mol	Joback Method
hf	-406.53	kJ/mol	Joback Method
hfus	21.01	kJ/mol	Joback Method
hvap	50.68	kJ/mol	Joback Method
log10ws	-3.08		Crippen Method
logp	2.930		Crippen Method
mcvol	172.220	ml/mol	McGowan Method
pc	2187.68	kPa	Joback Method
rinpol	1285.00		NIST Webbook
rinpol	1274.00		NIST Webbook
rinpol	1273.00		NIST Webbook
rinpol	1274.00		NIST Webbook
rinpol	1285.00		NIST Webbook
tb	557.02	K	Joback Method
tc	763.76	K	Joback Method
tf	280.34	K	Joback Method
vc	0.644	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	435.53	J/mol×K	557.02	Joback Method
cpg	455.28	J/mol×K	591.48	Joback Method
cpg	474.03	J/mol×K	625.93	Joback Method
cpg	491.78	J/mol×K	660.39	Joback Method
cpg	508.54	J/mol×K	694.85	Joback Method
cpg	524.33	J/mol×K	729.31	Joback Method
cpg	539.15	J/mol×K	763.76	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
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=C109010120&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/70-778-1/Isopulegol-acetate.pdf>

Generated by Cheméo on 2025-12-23 06:32:04.100089238 +0000 UTC m=+6219721.630129892.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.