

14-Hydroxymurolene

Inchi:	InChI=1S/C15H24O/c1-10(2)13-6-4-11(3)14-7-5-12(9-16)8-15(13)14/h4,8,10,13-16H,5-7
InchiKey:	DFIKBBWSGFHYMP-QLFBSQMISA-N
Formula:	C15H24O
SMILES:	CC1=CCC(C(C)C)C2C=C(CO)CCC12
Mol. weight [g/mol]:	220.35

Physical Properties

Property code	Value	Unit	Source
gf	42.21	kJ/mol	Joback Method
hf	-317.20	kJ/mol	Joback Method
hfus	25.78	kJ/mol	Joback Method
hvap	67.39	kJ/mol	Joback Method
log10ws	-3.89		Crippen Method
logp	3.554		Crippen Method
mcvol	197.760	ml/mol	McGowan Method
pc	2086.93	kPa	Joback Method
rinsol	1624.00		NIST Webbook
tb	668.51	K	Joback Method
tc	869.60	K	Joback Method
tf	348.75	K	Joback Method
vc	0.742	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	576.51	J/molxK	668.51	Joback Method
cpg	594.95	J/molxK	702.02	Joback Method
cpg	612.33	J/molxK	735.54	Joback Method
cpg	628.69	J/molxK	769.05	Joback Method
cpg	644.06	J/molxK	802.57	Joback Method
cpg	658.50	J/molxK	836.08	Joback Method
cpg	672.04	J/molxK	869.60	Joback Method
dvisc	0.0048013	Paxs	348.75	Joback Method
dvisc	0.0016816	Paxs	402.04	Joback Method

dvisc	0.0007529	Paxs	455.34	Joback Method
dvisc	0.0003989	Paxs	508.63	Joback Method
dvisc	0.0002384	Paxs	561.92	Joback Method
dvisc	0.0001558	Paxs	615.22	Joback Method
dvisc	0.0001089	Paxs	668.51	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=R202881&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
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
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/17-147-0/14-Hydroxymurolene.pdf>

Generated by Cheméo on 2024-04-26 05:44:49.167030528 +0000 UTC m=+16399538.087607841.

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