

(Z)-Myroxide

Other names:	(Z)-Miroxide cis-Myroxide
Inchi:	InChI=1S/C10H16O/c1-5-8(2)6-7-9-10(3,4)11-9/h5-6,9H,1,7H2,2-4H3/b8-6+
InchiKey:	LIMXJCIGROLRED-SOFGYWHQSA-N
Formula:	C10H16O
SMILES:	<chem>C=CC(C)=CCC1OC1(C)C</chem>
Mol. weight [g/mol]:	152.23
CAS:	94607-48-4

Physical Properties

Property code	Value	Unit	Source
gf	154.26	kJ/mol	Joback Method
hf	-81.17	kJ/mol	Joback Method
hfus	20.15	kJ/mol	Joback Method
hvap	40.19	kJ/mol	Joback Method
log10ws	-2.92		Crippen Method
logp	2.686		Crippen Method
mcvol	138.170	ml/mol	McGowan Method
pc	2640.67	kPa	Joback Method
rinpol	1137.00		NIST Webbook
rinpol	1125.00		NIST Webbook
rinpol	1134.00		NIST Webbook
rinpol	1114.00		NIST Webbook
rinpol	1137.00		NIST Webbook
rinpol	1131.00		NIST Webbook
rinpol	1137.00		NIST Webbook
rinpol	1127.00		NIST Webbook
rinpol	1132.00		NIST Webbook
rinpol	1131.00		NIST Webbook
ripol	1550.00		NIST Webbook
ripol	1493.00		NIST Webbook
ripol	1493.00		NIST Webbook
ripol	1550.00		NIST Webbook
tb	458.18	K	Joback Method
tc	657.47	K	Joback Method
tf	245.83	K	Joback Method
vc	0.532	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	302.19	J/mol×K	458.18	Joback Method
cpg	318.21	J/mol×K	491.40	Joback Method
cpg	333.07	J/mol×K	524.61	Joback Method
cpg	346.90	J/mol×K	557.83	Joback Method
cpg	359.80	J/mol×K	591.04	Joback Method
cpg	371.90	J/mol×K	624.26	Joback Method
cpg	383.32	J/mol×K	657.47	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C94607484&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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

Latest version available from:

<https://www.chemeo.com/cid/28-956-0/Z-Myroxide.pdf>

Generated by Cheméo on 2024-04-23 06:13:14.039132665 +0000 UTC m=+16142042.959709977.

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