

Oxirane, methyl-, (S)-

Other names:	(S)-(-)-Propylene oxide (S)-1,2-epoxypropane 1,2-Propylene oxide
Inchi:	InChI=1S/C3H6O/c1-3-2-4-3/h3H,2H2,1H3/t3-/m1/s1
InchiKey:	GOOHAUXETOMSMG-GSVUUGTGSA-N
Formula:	C3H6O
SMILES:	CC1CO1
Mol. weight [g/mol]:	58.08
CAS:	16088-62-3

Physical Properties

Property code	Value	Unit	Source
gf	-50.99	kJ/mol	Joback Method
hf	-164.45	kJ/mol	Joback Method
hfus	9.64	kJ/mol	Joback Method
hvap	26.70	kJ/mol	Joback Method
log10ws	-0.17		Crippen Method
logp	0.405		Crippen Method
mcvol	48.140	ml/mol	McGowan Method
pc	5213.15	kPa	Joback Method
tb	307.90 ± 2.00	K	NIST Webbook
tc	481.72	K	Joback Method
tf	168.08	K	Joback Method
vc	0.181	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	70.70	J/molxK	301.73	Joback Method
cpg	78.27	J/molxK	331.73	Joback Method
cpg	85.43	J/molxK	361.73	Joback Method
cpg	92.19	J/molxK	391.72	Joback Method
cpg	98.56	J/molxK	421.72	Joback Method
cpg	104.58	J/molxK	451.72	Joback Method

cpg	110.26	J/molxK	481.72	Joback Method
dvisc	0.0005257	Paxs	168.08	Joback Method
dvisc	0.0004254	Paxs	190.35	Joback Method
dvisc	0.0003599	Paxs	212.63	Joback Method
dvisc	0.0003143	Paxs	234.91	Joback Method
dvisc	0.0002810	Paxs	257.18	Joback Method
dvisc	0.0002557	Paxs	279.46	Joback Method
dvisc	0.0002360	Paxs	301.73	Joback Method

Sources

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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C16088623&Units=SI
The Yaws Handbook of Vapor Pressure:	https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure

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

Latest version available from:

<https://www.chemeo.com/cid/33-308-3/Oxirane-methyl-S.pdf>

Generated by Cheméo on 2022-12-03 21:42:33.30172937 +0000 UTC m=+100116.038595064.

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