

1,3-Oxathiolane, 2-ethyl-5-methyl, cis-

Inchi:	InChI=1S/C5H10OS/c1-4-3-7-5(2)6-4/h4-5H,3H2,1-2H3/t4-,5-/m1/s1
InchiKey:	HLLCFXLPRUGRGF-RFZPGFLSSA-N
Formula:	C6H12OS
SMILES:	CC1CSC(C)O1
Mol. weight [g/mol]:	132.22
CAS:	38384-66-6

Physical Properties

Property code	Value	Unit	Source
gf	-26.20	kJ/mol	Joback Method
hf	-193.13	kJ/mol	Joback Method
hfus	15.35	kJ/mol	Joback Method
hvap	36.99	kJ/mol	Joback Method
log10ws	-1.50		Crippen Method
logp	1.484		Crippen Method
mcvol	92.670	ml/mol	McGowan Method
pc	4077.71	kPa	Joback Method
tb	399.19	K	Joback Method
tc	613.97	K	Joback Method
tf	262.79	K	Joback Method
vc	0.323	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	172.65	J/molxK	399.19	Joback Method
cpg	185.37	J/molxK	434.99	Joback Method
cpg	197.47	J/molxK	470.78	Joback Method
cpg	208.94	J/molxK	506.58	Joback Method
cpg	219.82	J/molxK	542.38	Joback Method
cpg	230.11	J/molxK	578.17	Joback Method
cpg	239.83	J/molxK	613.97	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=C38384666&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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
hvac:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mccol:	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/17-548-5/1-3-Oxathiolane-2-ethyl-5-methyl-cis.pdf>

Generated by Cheméo on 2024-04-17 01:38:06.240541937 +0000 UTC m=+15607135.161119249.

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