

2-Isopropoxy-tetrahydro-furan

Other names:	Tetrahydrofuran, 2-(1-methylethyloxy)
Inchi:	InChI=1S/C7H14O2/c1-6(2)9-7-4-3-5-8-7/h6-7H,3-5H2,1-2H3
InchiKey:	OGAUHPGWQSCXSW-UHFFFAOYSA-N
Formula:	C7H14O2
SMILES:	CC(C)OC1CCCO1
Mol. weight [g/mol]:	130.18

Physical Properties

Property code	Value	Unit	Source
gf	-148.95	kJ/mol	Joback Method
hf	-396.83	kJ/mol	Joback Method
hfus	13.46	kJ/mol	Joback Method
hvap	37.97	kJ/mol	Joback Method
log10ws	-1.54		Crippen Method
logp	1.548		Crippen Method
mcvol	110.370	ml/mol	McGowan Method
pc	3348.98	kPa	Joback Method
rinpol	830.00		NIST Webbook
rinpol	830.00		NIST Webbook
tb	423.77	K	Joback Method
tc	622.67	K	Joback Method
tf	213.35	K	Joback Method
vc	0.402	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	229.36	J/molxK	423.77	Joback Method
cpg	244.23	J/molxK	456.92	Joback Method
cpg	258.44	J/molxK	490.07	Joback Method
cpg	272.01	J/molxK	523.22	Joback Method
cpg	284.95	J/molxK	556.37	Joback Method
cpg	297.27	J/molxK	589.52	Joback Method
cpg	308.97	J/molxK	622.67	Joback Method

dvisc	0.0059385	Paxs	213.35	Joback Method
dvisc	0.0025675	Paxs	248.42	Joback Method
dvisc	0.0013660	Paxs	283.49	Joback Method
dvisc	0.0008351	Paxs	318.56	Joback Method
dvisc	0.0005628	Paxs	353.63	Joback Method
dvisc	0.0004074	Paxs	388.70	Joback Method
dvisc	0.0003110	Paxs	423.77	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=R91070&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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
mvol:	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

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