Furan, tetrahydro-2,4-dimethyl-, cis-

Other names:	Furan, tetrahydro, 2,4-cis-dimethyl
	2,4-Dimethyltetrahydrofuran, (Z)-
Inchi:	InChI=1S/C6H12O/c1-5-3-6(2)7-4-5/h5-6H,3-4H2,1-2H3/t5-,6+/m0/s1
InchiKey:	QMGLMRPHOITLSN-NTSWFWBYSA-N
Formula:	C6H12O
SMILES:	CC1COC(C)C1
Mol. weight [g/mol]:	100.16
CAS:	39168-01-9

Physical Properties

Property code	Value	Unit	Source
gf	-57.64	kJ/mol	Joback Method
hf	-259.03	kJ/mol	Joback Method
hfus	14.28	kJ/mol	Joback Method
hvap	33.41	kJ/mol	Joback Method
log10ws	-1.19		Crippen Method
logp	1.431		Crippen Method
mcvol	90.410	ml/mol	McGowan Method
pc	3646.53	kPa	Joback Method
rinpol	1004.00		NIST Webbook
tb	374.24	K	Joback Method
tc	569.86	K	Joback Method
tf	190.61	K	Joback Method
vc	0.333	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	165.98	J/mol×K	374.24	Joback Method
cpg	179.78	J/mol×K	406.84	Joback Method
cpg	192.97	J/mol×K	439.45	Joback Method
cpg	205.56	J/mol×K	472.05	Joback Method
cpg	217.59	J/mol×K	504.65	Joback Method
cpg	229.04	J/mol×K	537.25	Joback Method

cpg	239.94	J/mol×K	569.86	Joback Method
dvisc	0.0023244	Pa×s	190.61	Joback Method
dvisc	0.0013336	Pa×s	221.22	Joback Method
dvisc	0.0008757	Pa×s	251.82	Joback Method
dvisc	0.0006300	Pa×s	282.43	Joback Method
dvisc	0.0004833	Pa×s	313.03	Joback Method
dvisc	0.0003887	Pa×s	343.63	Joback Method
dvisc	0.0003240	Pa×s	374.24	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
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=C39168019&Units=SI

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

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