

Tetrahydrofuran, 2-ethenyloxy

Other names:	2-Vinyloxy-tetrahydro-furan
Inchi:	InChI=1S/C6H10O2/c1-2-7-6-4-3-5-8-6/h2,6H,1,3-5H2
InchiKey:	ZNYVTSYLHMCSSI-UHFFFAOYSA-N
Formula:	C6H10O2
SMILES:	C=COC1CCCO1
Mol. weight [g/mol]:	114.14

Physical Properties

Property code	Value	Unit	Source
gf	-67.09	kJ/mol	Joback Method
hf	-245.48	kJ/mol	Joback Method
hfus	13.12	kJ/mol	Joback Method
hvap	35.46	kJ/mol	Joback Method
log10ws	-1.36		Crippen Method
logp	1.283		Crippen Method
mcvol	91.980	ml/mol	McGowan Method
pc	3916.03	kPa	Joback Method
rinpol	830.00		NIST Webbook
rinpol	830.00		NIST Webbook
rinpol	830.00		NIST Webbook
tb	398.01	K	Joback Method
tc	598.42	K	Joback Method
tf	215.32	K	Joback Method
vc	0.333	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	174.69	J/mol×K	398.01	Joback Method
cpg	187.21	J/mol×K	431.41	Joback Method
cpg	199.16	J/mol×K	464.81	Joback Method
cpg	210.54	J/mol×K	498.22	Joback Method
cpg	221.37	J/mol×K	531.62	Joback Method
cpg	231.65	J/mol×K	565.02	Joback Method

cpg	241.41	J/mol×K	598.42	Joback Method
dvisc	0.0031583	Pa×s	215.32	Joback Method
dvisc	0.0017063	Pa×s	245.77	Joback Method
dvisc	0.0010559	Pa×s	276.22	Joback Method
dvisc	0.0007187	Pa×s	306.66	Joback Method
dvisc	0.0005244	Pa×s	337.11	Joback Method
dvisc	0.0004032	Pa×s	367.56	Joback Method
dvisc	0.0003227	Pa×s	398.01	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R91137&Units=SI
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

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

Latest version available from:

<https://www.chemeo.com/cid/80-979-7/Tetrahydrofuran-2-ethenyloxy.pdf>

Generated by Cheméo on 2025-12-05 13:54:21.885991929 +0000 UTC m=+4691059.416032584.

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