

tert-Butyl glycidyl ether

Other names:	(tert-Butoxymethyl)oxirane t-Butyl glycidyl ether Oxirane, [(1,1-dimethylethoxy)methyl]- t-Bge 1,1-Dimethylethyl glycidyl ether Propane, 1-tert-butoxy-2,3-epoxy- ((1,1-Dimethylethoxy)methyl)oxirane Ageflex TBGE Glycidyl tert-butyl ether 2-(tert-Butoxymethyl)oxirane 2,3-Epoxypropyl-t-butyl ether Oxirane, 2-((1,1-dimethylethoxy)methyl)-
Inchi:	InChI=1S/C7H14O2/c1-7(2,3)9-5-6-4-8-6/h6H,4-5H2,1-3H3
InchiKey:	SFJRJUJEMVAZLM-UHFFFAOYSA-N
Formula:	C7H14O2
SMILES:	CC(C)(C)OCC1CO1
Mol. weight [g/mol]:	130.18
CAS:	7665-72-7

Physical Properties

Property code	Value	Unit	Source
chl	-4387.00 ± 1.00	kJ/mol	NIST Webbook
chl	-4385.00 ± 3.00	kJ/mol	NIST Webbook
chl	-4387.00 ± 1.00	kJ/mol	NIST Webbook
gf	-119.47	kJ/mol	Joback Method
hf	-319.70 ± 2.70	kJ/mol	NIST Webbook
hf	-319.80 ± 3.00	kJ/mol	NIST Webbook
hf	-318.80 ± 1.10	kJ/mol	NIST Webbook
hf	-322.00 ± 2.00	kJ/mol	NIST Webbook
hf	-318.80 ± 1.10	kJ/mol	NIST Webbook
hfl	-369.00 ± 1.00	kJ/mol	NIST Webbook
hfl	-370.00 ± 3.00	kJ/mol	NIST Webbook
hfl	-369.00 ± 1.00	kJ/mol	NIST Webbook
hfus	13.77	kJ/mol	Joback Method
hvac	50.22 ± 0.35	kJ/mol	NIST Webbook
log10ws	-1.05		Crippen Method
logp	1.200		Crippen Method

mvol	110.370	ml/mol	McGowan Method
pc	3202.78	kPa	Joback Method
tb	412.44	K	Joback Method
tc	603.87	K	Joback Method
tf	237.81	K	Joback Method
vc	0.412	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	233.77	J/mol×K	412.44	Joback Method
cpg	247.92	J/mol×K	444.34	Joback Method
cpg	261.29	J/mol×K	476.25	Joback Method
cpg	273.92	J/mol×K	508.15	Joback Method
cpg	285.83	J/mol×K	540.06	Joback Method
cpg	297.07	J/mol×K	571.96	Joback Method
cpg	307.66	J/mol×K	603.87	Joback Method
dvisc	0.0027913	Paxs	237.81	Joback Method
dvisc	0.0017301	Paxs	266.92	Joback Method
dvisc	0.0011781	Paxs	296.02	Joback Method
dvisc	0.0008593	Paxs	325.12	Joback Method
dvisc	0.0006602	Paxs	354.23	Joback Method
dvisc	0.0005279	Paxs	383.34	Joback Method
dvisc	0.0004357	Paxs	412.44	Joback Method

Sources

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=C7665727&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

chl: Standard liquid enthalpy of combustion

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfl:	Liquid phase 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

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