2-Butyl ethoxymethyl ether

Other names: Methane, 2-butoxy-ethoxy
Ethoxy-(2-butoxy)methane
Inchi: InChI=1S/C7H16O2/c1-4-7(3)9-6-8-5-2/h7H,4-6H2,1-3H3
InchiKey: WLCHNYQQRXSZER-UHFFFAOYSA-N
Formula: C7H16O2
SMILES: CCOCOC(C)CC
Mol. weight [g/mol]: 132.20

Physical Properties

Property code	Value	Unit	Source
gf	-204.38	kJ/mol	Joback Method
hf	-457.53	kJ/mol	Joback Method
hfus	12.74	kJ/mol	Joback Method
hvap	35.61	kJ/mol	Joback Method
log10ws	-1.53		Crippen Method
logp	1.796		Crippen Method
mcvol	121.230	ml/mol	McGowan Method
pc	2724.01	kPa	Joback Method
rinpol	794.00		NIST Webbook
rinpol	826.00		NIST Webbook
rinpol	826.00		NIST Webbook
rinpol	803.00		NIST Webbook
ripol	992.00		NIST Webbook
tb	403.96	K	Joback Method
tc	572.52	K	Joback Method
tf	198.11	K	Joback Method
vc	0.458	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	242.30	J/mol×K	403.96	Joback Method
cpg	253.90	J/mol×K	432.05	Joback Method
cpg	265.20	J/mol×K	460.15	Joback Method

cpg	276.21	J/mol×K	488.24	Joback Method
cpg	286.90	J/mol×K	516.33	Joback Method
cpg	297.29	J/mol×K	544.43	Joback Method
cpg	307.36	J/mol×K	572.52	Joback Method
dvisc	0.0047999	Pa×s	198.11	Joback Method
dvisc	0.0018890	Pa×s	232.42	Joback Method
dvisc	0.0009450	Pa×s	266.73	Joback Method
dvisc	0.0005536	Pa×s	301.04	Joback Method
dvisc	0.0003618	Pa×s	335.34	Joback Method
dvisc	0.0002559	Pa×s	369.65	Joback Method
dvisc	0.0001919	Pa×s	403.96	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R143574&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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