

Propyl tert-pentyl ether

Other names:	1,1-dimethylpropyl propyl ether 2-methyl-2-propoxybutane butane, 2-methyl-2-propoxy- tert-pentyl propyl ether
Inchi:	InChI=1S/C8H18O/c1-5-7-9-8(3,4)6-2/h5-7H2,1-4H3
InchiKey:	LFKBOUUQIPXNTJ-UHFFFAOYSA-N
Formula:	C8H18O
SMILES:	CCCOC(C)(C)CC
Mol. weight [g/mol]:	130.23

Physical Properties

Property code	Value	Unit	Source
gf	-85.68	kJ/mol	Joback Method
hf	-349.42	kJ/mol	Joback Method
hfus	10.25	kJ/mol	Joback Method
hvap	34.52	kJ/mol	Joback Method
log10ws	-2.37		Crippen Method
logp	2.602		Crippen Method
mcvol	129.450	ml/mol	McGowan Method
pc	2540.49	kPa	Joback Method
rinpol	812.00		NIST Webbook
tb	401.63	K	Joback Method
tc	576.09	K	Joback Method
tf	204.57	K	Joback Method
vc	0.490	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	335.39	J/molxK	576.09	Joback Method
cpg	273.09	J/molxK	430.71	Joback Method
cpg	286.65	J/molxK	459.78	Joback Method
cpg	299.64	J/molxK	488.86	Joback Method
cpg	312.08	J/molxK	517.94	Joback Method

cpg	323.99	J/molxK	547.02	Joback Method
cpg	258.96	J/molxK	401.63	Joback Method
dvisc	0.0073558	Paxs	204.57	Joback Method
dvisc	0.0028037	Paxs	237.41	Joback Method
dvisc	0.0013510	Paxs	270.26	Joback Method
dvisc	0.0007626	Paxs	303.10	Joback Method
dvisc	0.0004814	Paxs	335.94	Joback Method
dvisc	0.0003298	Paxs	368.79	Joback Method
dvisc	0.0002404	Paxs	401.63	Joback Method
pvap	0.40	kPa	274.20	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	0.44	kPa	276.20	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	0.53	kPa	278.10	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	0.61	kPa	280.20	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	0.72	kPa	282.20	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	0.79	kPa	284.20	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	0.80	kPa	285.10	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers

pvap	1.00	kPa	287.30	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	1.20	kPa	290.30	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	1.43	kPa	293.30	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	1.80	kPa	296.20	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	2.07	kPa	299.20	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers
pvap	2.33	kPa	302.20	Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers

Sources

- Thermochemistry of Branched Ethers: Experimental Study of Chemical Equilibria in the Reacting System of tert-Amyl Alkyl Ether Synthesis: McGowan Method:** <https://www.doi.org/10.1021/je034172y>
- Joback Method:** https://en.wikipedia.org/wiki/Joback_method
- Springer Method:** <http://link.springer.com/article/10.1007/BF02311772>
- NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=R559865&Units=SI>
- Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
- Crippen Method:** https://www.chemeo.com/doc/models/crippen_log10ws
- Determination of Ambient Temperature Vapor Pressures and Vaporization Enthalpies of Branched Ethers:** <https://www.doi.org/10.1021/je0255980>

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
pvap:	Vapor pressure
rinpolar:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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