

(5-ETHYLCYCLOPENT-1-ENYL)METHANOL

Inchi:

InChI=1S/C8H14O/c1-2-7-4-3-5-8(7)6-9/h5,7,9H,2-4,6H2,1H3

InchiKey:

PWPHPRWZLTZIPI-UHFFFAOYSA-N

Formula:

C8H14O

SMILES:

CCC1CCC=C1CO

Mol. weight [g/mol]:

126.20

Physical Properties

Property code	Value	Unit	Source
hfus	15.33	kJ/mol	Joback Method
logp	1.725		Crippen Method
mcvol	114.290	ml/mol	McGowan Method
rinpol	1073.00		NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	258.78	J/mol×K	494.04	Joback Method
cpg	271.32	J/mol×K	525.16	Joback Method
cpg	283.25	J/mol×K	556.28	Joback Method
cpg	294.60	J/mol×K	587.40	Joback Method
cpg	305.39	J/mol×K	618.52	Joback Method
cpg	315.63	J/mol×K	649.64	Joback Method
cpg	325.35	J/mol×K	680.76	Joback Method
dvisc	0.0172801	Pa×s	264.92	Joback Method
dvisc	0.0051460	Pa×s	303.11	Joback Method
dvisc	0.0020097	Pa×s	341.29	Joback Method
dvisc	0.0009483	Pa×s	379.48	Joback Method
dvisc	0.0005134	Pa×s	417.67	Joback Method
dvisc	0.0003080	Pa×s	455.85	Joback Method
dvisc	0.0002000	Pa×s	494.04	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C36431591&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
hfus:	Enthalpy of fusion at standard conditions
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpola:	Non-polar retention indices

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