8-hexadecenal, E

Other names: (E)-8-Hexadecenal
Inchi: InChI=1S/C16H30O/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-16-17/h8-9,16H,2-7,10-15H2,
InchiKey: CDLTZHMYSOPNNK-CMDGGOBGSA-N
Formula: C16H30O
SMILES: CCCCCCCC=CCCCCCCC=O
Mol. weight [g/mol]: 238.41

Physical Properties

Property code	Value	Unit	Source
gf	64.54	kJ/mol	Joback Method
hf	-341.93	kJ/mol	Joback Method
hfus	39.69	kJ/mol	Joback Method
hvap	57.89	kJ/mol	Joback Method
log10ws	-5.65		Crippen Method
logp	5.443		Crippen Method
mcvol	233.570	ml/mol	McGowan Method
pc	1447.94	kPa	Joback Method
rinpol	1803.00		NIST Webbook
rinpol	1803.00		NIST Webbook
ripol	2144.00		NIST Webbook
ripol	2144.00		NIST Webbook
tb	618.30	K	Joback Method
tc	787.48	K	Joback Method
tf	307.00	K	Joback Method
vc	0.928	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	624.39	J/mol×K	618.30	Joback Method
cpg	641.87	J/mol×K	646.50	Joback Method
cpg	658.57	J/mol×K	674.69	Joback Method
cpg	674.53	J/mol×K	702.89	Joback Method
cpg	689.77	J/mol×K	731.09	Joback Method

cpg	704.32	J/mol×K	759.29	Joback Method
cpg	718.22	J/mol×K	787.48	Joback Method
dvisc	0.0038359	Pa×s	307.00	Joback Method
dvisc	0.0015057	Pa×s	358.88	Joback Method
dvisc	0.0007485	Pa×s	410.77	Joback Method
dvisc	0.0004352	Pa×s	462.65	Joback Method
dvisc	0.0002823	Pa×s	514.53	Joback Method
dvisc	0.0001983	Pa×s	566.42	Joback Method
dvisc	0.0001477	Pa×s	618.30	Joback Method

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R265451&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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