cis-3-Hexenoic acid

Other names:	(Z)-3-hexenoic acid
	(Z)-hex-2-enoic acid
Inchi:	InChI=1S/C6H10O2/c1-2-3-4-5-6(7)8/h3-4H,2,5H2,1H3,(H,7,8)/b4-3-
InchiKey:	XXHDAWYDNSXJQM-ARJAWSKDSA-N
Formula:	C6H10O2
SMILES:	CCC=CCC(=O)O
Mol. weight [g/mol]:	114.14
CAS:	1775-43-5

Physical Properties

Property code	Value	Unit	Source
gf	-185.88	kJ/mol	Joback Method
hf	-314.76	kJ/mol	Joback Method
hfus	17.19	kJ/mol	Joback Method
hvap	52.33	kJ/mol	Joback Method
log10ws	-1.29		Crippen Method
logp	1.427		Crippen Method
mcvol	98.540	ml/mol	McGowan Method
pc	3970.51	kPa	Joback Method
rinpol	1101.00		NIST Webbook
ripol	1945.00		NIST Webbook
ripol	1922.00		NIST Webbook
tb	486.89	K	Joback Method
tc	665.94	K	Joback Method
tf	263.05	K	Joback Method
vc	0.377	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	206.08	J/mol×K	486.89	Joback Method
cpg	214.50	J/mol×K	516.73	Joback Method
cpg	222.51	J/mol×K	546.57	Joback Method
cpg	230.13	J/mol×K	576.41	Joback Method

cpg	237.37	J/mol×K	606.26	Joback Method
cpg	244.26	J/mol×K	636.10	Joback Method
cpg	250.80	J/mol×K	665.94	Joback Method
dvisc	0.0232383	Pa×s	263.05	Joback Method
dvisc	0.0061439	Pa×s	300.36	Joback Method
dvisc	0.0021795	Pa×s	337.66	Joback Method
dvisc	0.0009502	Pa×s	374.97	Joback Method
dvisc	0.0004814	Pa×s	412.28	Joback Method
dvisc	0.0002731	Pa×s	449.58	Joback Method
dvisc	0.0001689	Pa×s	486.89	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=C1775435&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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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