

2-Hexene, 1,1-diethoxy-

Other names:	1,1-Diethoxy-2-hexene 1,1-diethoxyhex-2-ene
Inchi:	InChI=1S/C10H20O2/c1-4-7-8-9-10(11-5-2)12-6-3/h8-10H,4-7H2,1-3H3/b9-8+
InchiKey:	WMQKYHTZGYIHHD-CMDGGGOBGSA-N
Formula:	C10H20O2
SMILES:	CCCC=CC(OCC)OCC
Mol. weight [g/mol]:	172.26
CAS:	54306-00-2

Physical Properties

Property code	Value	Unit	Source
gf	-98.90	kJ/mol	Joback Method
hf	-402.23	kJ/mol	Joback Method
hfus	20.71	kJ/mol	Joback Method
hvap	42.24	kJ/mol	Joback Method
log10ws	-2.64		Crippen Method
logp	2.742		Crippen Method
mvol	159.200	ml/mol	McGowan Method
pc	2159.31	kPa	Joback Method
tb	476.76	K	Joback Method
tc	650.07	K	Joback Method
tf	226.84	K	Joback Method
vc	0.606	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	353.11	J/mol×K	476.76	Joback Method
cpg	367.77	J/mol×K	505.64	Joback Method
cpg	381.91	J/mol×K	534.53	Joback Method
cpg	395.52	J/mol×K	563.41	Joback Method
cpg	408.61	J/mol×K	592.30	Joback Method
cpg	421.20	J/mol×K	621.18	Joback Method
cpg	433.28	J/mol×K	650.07	Joback Method

dvisc	0.0044403	Paxs	226.84	Joback Method
dvisc	0.0015806	Paxs	268.49	Joback Method
dvisc	0.0007426	Paxs	310.15	Joback Method
dvisc	0.0004172	Paxs	351.80	Joback Method
dvisc	0.0002648	Paxs	393.45	Joback Method
dvisc	0.0001834	Paxs	435.11	Joback Method
dvisc	0.0001354	Paxs	476.76	Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	369.70	K	4.70	NIST Webbook

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C54306002&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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
mccvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tbrp:	Boiling point at reduced pressure

tc: Critical Temperature
tf: Normal melting (fusion) point
vc: Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/46-517-7/2-Hexene-1-1-diethoxy.pdf>

Generated by Cheméo on 2024-04-29 00:17:43.476350766 +0000 UTC m=+16639112.396928087.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.