Undecanol-3

Other names:	3-Undecanol
	Undecan-3-ol
Inchi:	InChI=1S/C11H24O/c1-3-5-6-7-8-9-10-11(12)4-2/h11-12H,3-10H2,1-2H3
InchiKey:	HCARCYFXWDRVBZ-UHFFFAOYSA-N
Formula:	C11H24O
SMILES:	CCCCCCCCC(O)CC
Mol. weight [g/mol]:	172.31
CAS:	6929-08-4

Physical Properties

Property code	Value	Unit	Source
gf	-97.52	kJ/mol	Joback Method
hf	-427.88	kJ/mol	Joback Method
hfus	24.81	kJ/mol	Joback Method
hvap	56.37	kJ/mol	Joback Method
log10ws	-3.80		Crippen Method
logp	3.508		Crippen Method
mcvol	171.720	ml/mol	McGowan Method
pc	2115.83	kPa	Joback Method
rinpol	1400.00		NIST Webbook
rinpol	1308.00		NIST Webbook
rinpol	1308.00		NIST Webbook
rinpol	1400.00		NIST Webbook
ripol	1673.00		NIST Webbook
ripol	1665.00		NIST Webbook
ripol	1700.00		NIST Webbook
ripol	1665.00		NIST Webbook
tb	502.15 ± 4.00	K	NIST Webbook
tc	703.49	K	Joback Method
tf	259.55	K	Joback Method
vc	0.664	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	428.91	J/mol×K	542.82	Joback Method
cpg	443.16	J/mol×K	569.60	Joback Method
cpg	456.84	J/mol×K	596.38	Joback Method
cpg	469.99	J/mol×K	623.16	Joback Method
cpg	482.61	J/mol×K	649.93	Joback Method
cpg	494.72	J/mol×K	676.71	Joback Method
cpg	506.33	J/mol×K	703.49	Joback Method
dvisc	0.0500348	Pa×s	259.55	Joback Method
dvisc	0.0079438	Pa×s	306.76	Joback Method
dvisc	0.0020606	Pa×s	353.97	Joback Method
dvisc	0.0007343	Pa×s	401.18	Joback Method
dvisc	0.0003252	Pa×s	448.40	Joback Method
dvisc	0.0001682	Pa×s	495.61	Joback Method
dvisc	0.0000976	Pa×s	542.82	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	ln(Pvp) = A + B/(T + C)
Coeff. A	1.60089e+01
Coeff. B	-4.84174e+03
Coeff. C	-8.07480e+01
Temperature range (K), min.	388.72
Temperature range (K), max.	533.35

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=C6929084&Units=SI

The Yaws Handbook of Vapor Pressure: https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure

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
pvap:	Vapor 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