

2-Tridecen-1-ol, (E)-

Other names:

(2E)-2-Tridecen-1-ol

(Z)-tridec-2-en-1-ol

E-2-Tridecen-1-ol

trans-2-Tridecen-1-ol

Inchi:

InChI=1S/C13H26O/c1-2-3-4-5-6-7-8-9-10-11-12-13-14/h11-12,14H,2-10,13H2,1H3/b12-

InchiKey:

VPYJHNADOJDSGU-VAWYXSNFSA-N

Formula:

C13H26O

SMILES:

CCCCCCCCC=CCO

Mol. weight [g/mol]:

198.34

CAS:

74962-98-4

Physical Properties

Property code	Value	Unit	Source
gf	1.98	kJ/mol	Joback Method
hf	-346.66	kJ/mol	Joback Method
hfus	33.72	kJ/mol	Joback Method
hvap	61.17	kJ/mol	Joback Method
log10ws	-4.38		Crippen Method
logp	4.066		Crippen Method
mcvol	195.600	ml/mol	McGowan Method
pc	1854.71	kPa	Joback Method
rinpol	1570.00		NIST Webbook
rinpol	1570.00		NIST Webbook
tb	593.18	K	Joback Method
tc	756.04	K	Joback Method
tf	292.01	K	Joback Method
vc	0.762	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	510.30	J/mol×K	593.18	Joback Method
cpg	525.14	J/mol×K	620.32	Joback Method
cpg	539.35	J/mol×K	647.47	Joback Method

cpg	552.96	J/mol×K	674.61	Joback Method
cpg	566.00	J/mol×K	701.75	Joback Method
cpg	578.48	J/mol×K	728.90	Joback Method
cpg	590.43	J/mol×K	756.04	Joback Method
dvisc	0.0165809	Paxs	292.01	Joback Method
dvisc	0.0032726	Paxs	342.20	Joback Method
dvisc	0.0009783	Paxs	392.40	Joback Method
dvisc	0.0003846	Paxs	442.60	Joback Method
dvisc	0.0001829	Paxs	492.79	Joback Method
dvisc	0.0000998	Paxs	542.99	Joback Method
dvisc	0.0000603	Paxs	593.18	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.45395e+01
Coeff. B	-4.74446e+03
Coeff. C	-9.56210e+01
Temperature range (K), min.	428.52
Temperature range (K), max.	609.76

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C74962984&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

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
rinpolar:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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