

Naphthalene, 1,1'-undecylidenebis[decahydro-

Other names:	1,1-Di(1'-decahydronaphthyl)undecane 1,1-Di-(decahydro-1-naphthyl)undecane 1,1-bis(decahydro-1-naphthyl)undecane Naphthalene, 1,1'-undecylidenebis*decahydro- Undecane, 1,1-bis(decahydro-1-naphthyl)-
Inchi:	InChI=1S/C31H56/c1-2-3-4-5-6-7-8-9-22-31(29-23-14-18-25-16-10-12-20-27(25)29)30-2
InchiKey:	BMIYKWRBLFKRLB-UHFFFAOYSA-N
Formula:	C31H56
SMILES:	CCCCCCCCC(C1CCCC2CCCC21)C1CCCC2CCCC21
Mol. weight [g/mol]:	428.78
CAS:	55373-96-1

Physical Properties

Property code	Value	Unit	Source
gf	338.48	kJ/mol	Joback Method
hf	-487.21	kJ/mol	Joback Method
hfus	50.40	kJ/mol	Joback Method
hvap	84.62	kJ/mol	Joback Method
log10ws	-10.69		Crippen Method
logp	10.346		Crippen Method
mcvol	404.210	ml/mol	McGowan Method
pc	801.60	kPa	Joback Method
tb	960.02	K	Joback Method
tc	1182.16	K	Joback Method
tf	459.25	K	Joback Method
vc	1.528	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1668.72	J/molxK	1182.16	Joback Method
cpg	1650.67	J/molxK	1145.14	Joback Method
cpg	1631.28	J/molxK	1108.12	Joback Method
cpg	1610.41	J/molxK	1071.09	Joback Method

cpg	1587.93	J/mol×K	1034.07	Joback Method
cpg	1563.69	J/mol×K	997.04	Joback Method
cpg	1537.54	J/mol×K	960.02	Joback Method
dvisc	0.0028215	Paxs	459.25	Joback Method
dvisc	0.0001815	Paxs	960.02	Joback Method
dvisc	0.0002307	Paxs	876.56	Joback Method
dvisc	0.0003083	Paxs	793.10	Joback Method
dvisc	0.0004411	Paxs	709.63	Joback Method
dvisc	0.0006943	Paxs	626.17	Joback Method
dvisc	0.0012566	Paxs	542.71	Joback Method
hvapt	110.50	kJ/mol	543.00	NIST Webbook

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.73251e+01
Coeff. B	-9.55167e+03
Temperature range (K), min.	560.63
Temperature range (K), max.	795.07

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C55373961&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/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
h vap:	Enthalpy of vaporization at standard conditions
hvapt:	Enthalpy of vaporization at a given temperature
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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