

Heptadecane, 4-propyl-

Other names:	4-Propylheptadecane 4-n-Propylheptadecane
Inchi:	InChI=1S/C20H42/c1-4-7-8-9-10-11-12-13-14-15-16-19-20(17-5-2)18-6-3/h20H,4-19H2,1
InchiKey:	KWOQMARCKSGDGU-UHFFFAOYSA-N
Formula:	C20H42
SMILES:	CCCCCCCCCCCC(CCC)CCC
Mol. weight [g/mol]:	282.55
CAS:	55044-10-5

Physical Properties

Property code	Value	Unit	Source
gf	115.08	kJ/mol	Joback Method
hf	-461.41	kJ/mol	Joback Method
hfus	44.03	kJ/mol	Joback Method
hvap	59.73	kJ/mol	Joback Method
log10ws	-7.95		Crippen Method
logp	7.904		Crippen Method
mvol	292.660	ml/mol	McGowan Method
pc	1020.73	kPa	Joback Method
tb	656.56	K	Joback Method
tc	818.58	K	Joback Method
tf	229.70 ± 10.00	K	NIST Webbook
tf	270.70 ± 10.00	K	NIST Webbook
tf	270.65	K	NIST Webbook
vc	1.149	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	955.21	J/mol×K	818.58	Joback Method
cpg	938.17	J/mol×K	791.58	Joback Method
cpg	920.35	J/mol×K	764.58	Joback Method
cpg	901.73	J/mol×K	737.57	Joback Method
cpg	882.29	J/mol×K	710.57	Joback Method

cpg	862.00	J/mol×K	683.56	Joback Method
cpg	840.82	J/mol×K	656.56	Joback Method
dvisc	0.0051032	Paxs	300.16	Joback Method
dvisc	0.0000902	Paxs	656.56	Joback Method
dvisc	0.0001265	Paxs	597.16	Joback Method
dvisc	0.0001911	Paxs	537.76	Joback Method
dvisc	0.0003200	Paxs	478.36	Joback Method
dvisc	0.0006199	Paxs	418.96	Joback Method
dvisc	0.0014944	Paxs	359.56	Joback Method
hvapt	79.20	kJ/mol	442.00	NIST Webbook

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.42676e+01
Coeff. B	-4.77175e+03
Coeff. C	-1.02239e+02
Temperature range (K), min.	443.56
Temperature range (K), max.	635.03

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C55044105&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/ci990307I
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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