

Pentalene, octahydro-

Other names:	Bicyclo[3.3.0]octane Octahydropentalene
Inchi:	InChI=1S/C8H14/c1-3-7-5-2-6-8(7)4-1/h7-8H,1-6H2
InchiKey:	AEBWATHAIVJLTA-UHFFFAOYSA-N
Formula:	C8H14
SMILES:	C1CC2CCCC2C1
Mol. weight [g/mol]:	110.20
CAS:	694-72-4

Physical Properties

Property code	Value	Unit	Source
gf	113.78	kJ/mol	Joback Method
hf	-75.17	kJ/mol	Joback Method
hfus	8.55	kJ/mol	Joback Method
hvap	33.57	kJ/mol	Joback Method
log10ws	-2.48		Crippen Method
logp	2.587		Crippen Method
mvol	101.860	ml/mol	McGowan Method
pc	3560.02	kPa	Joback Method
rinpol	870.00		NIST Webbook
rinpol	870.00		NIST Webbook
tb	404.46	K	Joback Method
tc	614.38	K	Joback Method
tf	208.76	K	Joback Method
vc	0.382	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	198.70	J/mol×K	404.46	Joback Method
cpg	279.71	J/mol×K	579.39	Joback Method
cpg	265.55	J/mol×K	544.41	Joback Method
cpg	250.43	J/mol×K	509.42	Joback Method
cpg	234.29	J/mol×K	474.43	Joback Method

cpg	217.06	J/molxK	439.45	Joback Method
cpg	292.95	J/molxK	614.38	Joback Method
dvisc	0.0004754	Paxs	404.46	Joback Method
dvisc	0.0005251	Paxs	371.84	Joback Method
dvisc	0.0005912	Paxs	339.23	Joback Method
dvisc	0.0006827	Paxs	306.61	Joback Method
dvisc	0.0008158	Paxs	273.99	Joback Method
dvisc	0.0010230	Paxs	241.38	Joback Method
dvisc	0.0013767	Paxs	208.76	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C694724&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
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
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/46-148-7/Pentalene-octahydro.pdf>

Generated by Cheméo on 2024-04-25 21:00:19.200116932 +0000 UTC m=+16368068.120694244.

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