

2-phenyl-adamantane

Inchi:	InChI=1S/C16H20/c1-2-4-13(5-3-1)16-14-7-11-6-12(9-14)10-15(16)8-11/h1-5,11-12,14-1
InchiKey:	MWQZYLIXDPNHFG-PMONVTPFSA-N
Formula:	C16H20
SMILES:	<chem>c1ccc(C2C3CC4CC(C3)CC2C4)cc1</chem>
Mol. weight [g/mol]:	212.33

Physical Properties

Property code	Value	Unit	Source
gf	350.98	kJ/mol	Joback Method
hf	34.52	kJ/mol	Joback Method
hfus	25.68	kJ/mol	Joback Method
hvap	52.78	kJ/mol	Joback Method
log10ws	-4.30		Crippen Method
logp	4.226		Crippen Method
mcvol	179.960	ml/mol	McGowan Method
pc	2324.78	kPa	Joback Method
rinsol	1796.00		NIST Webbook
tb	607.31	K	Joback Method
tc	847.22	K	Joback Method
tf	338.32	K	Joback Method
vc	0.684	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	508.71	J/molxK	607.31	Joback Method
cpg	532.45	J/molxK	647.30	Joback Method
cpg	554.36	J/molxK	687.28	Joback Method
cpg	574.60	J/molxK	727.27	Joback Method
cpg	593.33	J/molxK	767.25	Joback Method
cpg	610.70	J/molxK	807.24	Joback Method
cpg	626.88	J/molxK	847.22	Joback Method
dvisc	0.0022093	Paxs	338.32	Joback Method
dvisc	0.0021934	Paxs	383.15	Joback Method

dvisc	0.0021809	Paxs	427.98	Joback Method
dvisc	0.0021708	Paxs	472.81	Joback Method
dvisc	0.0021625	Paxs	517.65	Joback Method
dvisc	0.0021556	Paxs	562.48	Joback Method
dvisc	0.0021497	Paxs	607.31	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R202227&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

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

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