

Indane, 1-(2-phenylethyl)

Inchi:	InChI=1S/C17H18/c1-2-6-14(7-3-1)10-11-16-13-12-15-8-4-5-9-17(15)16/h1-9,16H,10-13
InchiKey:	HHCVWDDVNPVAEC-UHFFFAOYSA-N
Formula:	C17H18
SMILES:	<chem>c1ccc(CCC2CCc3ccccc32)cc1</chem>
Mol. weight [g/mol]:	222.32

Physical Properties

Property code	Value	Unit	Source
gf	368.20	kJ/mol	Joback Method
hf	140.18	kJ/mol	Joback Method
hfus	25.61	kJ/mol	Joback Method
hvap	58.56	kJ/mol	Joback Method
log10ws	-4.97		Crippen Method
logp	4.349		Crippen Method
mcvol	192.010	ml/mol	McGowan Method
pc	2327.03	kPa	Joback Method
rinqol	1768.00		NIST Webbook
tb	653.44	K	Joback Method
tc	895.97	K	Joback Method
tf	364.65	K	Joback Method
vc	0.729	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	509.64	J/molxK	653.44	Joback Method
cpg	529.00	J/molxK	693.86	Joback Method
cpg	546.87	J/molxK	734.28	Joback Method
cpg	563.37	J/molxK	774.71	Joback Method
cpg	578.62	J/molxK	815.13	Joback Method
cpg	592.76	J/molxK	855.55	Joback Method
cpg	605.90	J/molxK	895.97	Joback Method
dvisc	0.0019234	Paxs	364.65	Joback Method
dvisc	0.0012250	Paxs	412.78	Joback Method

dvisc	0.0008573	Paxs	460.91	Joback Method
dvisc	0.0006418	Paxs	509.04	Joback Method
dvisc	0.0005052	Paxs	557.18	Joback Method
dvisc	0.0004130	Paxs	605.31	Joback Method
dvisc	0.0003479	Paxs	653.44	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R72995&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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