

P-phenylphenetole

Inchi:	InChI=1S/C14H14O/c1-2-15-14-10-8-13(9-11-14)12-6-4-3-5-7-12/h3-11H,2H2,1H3
InchiKey:	QRYTXFUBAZNNBP-UHFFFAOYSA-N
Formula:	C14H14O
SMILES:	CCOc1ccc(-c2ccccc2)cc1
Mol. weight [g/mol]:	198.26
CAS:	613-40-1

Physical Properties

Property code	Value	Unit	Source
gf	177.19	kJ/mol	Joback Method
hf	-2.92	kJ/mol	Joback Method
hfus	20.90	kJ/mol	Joback Method
hvap	54.38	kJ/mol	Joback Method
log10ws	-4.61		Crippen Method
logp	3.752		Crippen Method
mcvol	166.470	ml/mol	McGowan Method
pc	2679.08	kPa	Joback Method
tb	600.48	K	Joback Method
tc	837.55	K	Joback Method
tf	335.13	K	Joback Method
vc	0.622	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	397.75	J/molxK	600.48	Joback Method
cpg	414.65	J/molxK	639.99	Joback Method
cpg	430.37	J/molxK	679.50	Joback Method
cpg	444.95	J/molxK	719.01	Joback Method
cpg	458.45	J/molxK	758.53	Joback Method
cpg	470.90	J/molxK	798.04	Joback Method
cpg	482.36	J/molxK	837.55	Joback Method
dvisc	0.0014623	Paxs	335.13	Joback Method
dvisc	0.0007910	Paxs	379.36	Joback Method

dvisc	0.0004865	Paxs	423.58	Joback Method
dvisc	0.0003280	Paxs	467.81	Joback Method
dvisc	0.0002367	Paxs	512.03	Joback Method
dvisc	0.0001799	Paxs	556.25	Joback Method
dvisc	0.0001424	Paxs	600.48	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C613401&Units=SI

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
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/49-471-5/P-phenylphenetole.pdf>

Generated by Cheméo on 2024-04-19 17:55:35.573672454 +0000 UTC m=+15838584.494249764.

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