

2-Isopropylbicyclohexyl

Inchi:	InChI=1S/C15H28/c1-12(2)14-10-6-7-11-15(14)13-8-4-3-5-9-13/h12-15H,3-11H2,1-2H3
InchiKey:	XWULVJMSQHXLOJ-UHFFFAOYSA-N
Formula:	C15H28
SMILES:	CC(C)C1CCCCC1C1CCCCC1
Mol. weight [g/mol]:	208.38
CAS:	66374-73-0

Physical Properties

Property code	Value	Unit	Source
gf	114.17	kJ/mol	Joback Method
hf	-269.91	kJ/mol	Joback Method
hfus	15.82	kJ/mol	Joback Method
hvap	49.14	kJ/mol	Joback Method
log10ws	-4.92		Crippen Method
logp	5.029		Crippen Method
mcvol	200.490	ml/mol	McGowan Method
pc	1935.54	kPa	Joback Method
tb	550.33 ± 0.40	K	NIST Webbook
tb	556.21 ± 0.40	K	NIST Webbook
tc	800.93	K	Joback Method
tf	254.33	K	Joback Method
vc	0.735	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	688.24	J/mol×K	800.93	Joback Method
cpg	668.39	J/mol×K	763.54	Joback Method
cpg	647.02	J/mol×K	726.15	Joback Method
cpg	624.09	J/mol×K	688.76	Joback Method
cpg	599.56	J/mol×K	651.37	Joback Method
cpg	573.36	J/mol×K	613.98	Joback Method
cpg	545.45	J/mol×K	576.59	Joback Method
cpl	418.40	J/mol×K	313.00	NIST Webbook

dvisc	0.0002820	Paxs	522.88	Joback Method
dvisc	0.0004147	Paxs	469.17	Joback Method
dvisc	0.0006737	Paxs	415.46	Joback Method
dvisc	0.0012644	Paxs	361.75	Joback Method
dvisc	0.0029553	Paxs	308.04	Joback Method
dvisc	0.0098872	Paxs	254.33	Joback Method
dvisc	0.0002060	Paxs	576.59	Joback Method

Sources

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=C66374730&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.com/doc/models/crippen_log10ws

Legend

cpg:	Ideal gas heat capacity
cpl:	Liquid phase 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.cheméo.com/cid/23-960-0/2-Isopropylbicyclohexyl.pdf>

Generated by Cheméo on 2024-04-27 03:30:25.469540394 +0000 UTC m=+16477874.390117706.

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