

Cycloheptane, 1-isopropyl-

Inchi:	InChI=1S/C10H18/c1-9(2)10-7-5-3-4-6-8-10/h7,9H,3-6,8H2,1-2H3
InchiKey:	KTLJXRGLODFJIG-UHFFFAOYSA-N
Formula:	C10H18
SMILES:	CC(C)C1=CCCCC1
Mol. weight [g/mol]:	138.25
CAS:	17257-36-2

Physical Properties

Property code	Value	Unit	Source
gf	71.27	kJ/mol	Joback Method
hf	-140.20	kJ/mol	Joback Method
hfus	7.63	kJ/mol	Joback Method
hvap	39.33	kJ/mol	Joback Method
log10ws	-3.52		Crippen Method
logp	3.533		Crippen Method
mcvol	136.600	ml/mol	McGowan Method
pc	2811.36	kPa	Joback Method
tb	460.39	K	Joback Method
tc	673.79	K	Joback Method
tf	208.84	K	Joback Method
vc	0.501	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	286.32	J/molxK	460.39	Joback Method
cpg	305.44	J/molxK	495.96	Joback Method
cpg	323.57	J/molxK	531.52	Joback Method
cpg	340.73	J/molxK	567.09	Joback Method
cpg	356.95	J/molxK	602.65	Joback Method
cpg	372.25	J/molxK	638.22	Joback Method
cpg	386.66	J/molxK	673.79	Joback Method
dvisc	0.0177954	Paxs	208.84	Joback Method
dvisc	0.0044613	Paxs	250.76	Joback Method

dvisc	0.0016624	Paxs	292.69	Joback Method
dvisc	0.0007933	Paxs	334.62	Joback Method
dvisc	0.0004464	Paxs	376.54	Joback Method
dvisc	0.0002819	Paxs	418.47	Joback Method
dvisc	0.0001935	Paxs	460.39	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=C17257362&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/26-785-2/Cycloheptane-1-isopropyl.pdf>

Generated by Cheméo on 2024-04-23 11:36:00.232062928 +0000 UTC m=+16161409.152640241.

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