

(7,8-Methylene)octyl-cyclopropane

Inchi:	InChI=1S/C12H22/c1(3-5-11-7-8-11)2-4-6-12-9-10-12/h11-12H,1-10H2
InchiKey:	AKQNFNHJHSBCOS-UHFFFAOYSA-N
Formula:	C12H22
SMILES:	C(CCCC1CC1)CCC1CC1
Mol. weight [g/mol]:	166.30

Physical Properties

Property code	Value	Unit	Source
gf	171.66	kJ/mol	Joback Method
hf	-145.41	kJ/mol	Joback Method
hfus	23.11	kJ/mol	Joback Method
hvap	42.13	kJ/mol	Joback Method
log10ws	-4.15		Crippen Method
logp	4.147		Crippen Method
mcvol	158.220	ml/mol	McGowan Method
pc	2246.13	kPa	Joback Method
tb	487.44	K	Joback Method
tc	673.29	K	Joback Method
tf	260.88	K	Joback Method
vc	0.622	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	380.29	J/molxK	487.44	Joback Method
cpg	399.52	J/molxK	518.42	Joback Method
cpg	417.67	J/molxK	549.39	Joback Method
cpg	434.82	J/molxK	580.37	Joback Method
cpg	451.02	J/molxK	611.34	Joback Method
cpg	466.33	J/molxK	642.32	Joback Method
cpg	480.81	J/molxK	673.29	Joback Method
dvisc	0.0016523	Paxs	260.88	Joback Method
dvisc	0.0013554	Paxs	298.64	Joback Method
dvisc	0.0011624	Paxs	336.40	Joback Method

dvisc	0.0010282	Paxs	374.16	Joback Method
dvisc	0.0009303	Paxs	411.92	Joback Method
dvisc	0.0008559	Paxs	449.68	Joback Method
dvisc	0.0007977	Paxs	487.44	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=R137732&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/79-427-1/7-8-Methylene-octyl-cyclopropane.pdf>

Generated by Cheméo on 2024-04-30 09:02:04.48885569 +0000 UTC m=+16756973.409433002.

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