

5,8-Endomethylene-5,8-dihydro-1,4-naphthohydro

Inchi:	InChI=1S/C11H10O2/c12-8-3-4-9(13)11-7-2-1-6(5-7)10(8)11/h1-4,6-7,12-13H,5H2
InchiKey:	PQZMLJBZPCMYDX-UHFFFAOYSA-N
Formula:	C11H10O2
SMILES:	Oc1ccc(O)c2c1C1C=CC2C1
Mol. weight [g/mol]:	174.20
CAS:	3090-47-9

Physical Properties

Property code	Value	Unit	Source
gf	-1.16	kJ/mol	Joback Method
hf	-190.39	kJ/mol	Joback Method
hfus	29.06	kJ/mol	Joback Method
hvap	68.99	kJ/mol	Joback Method
log10ws	-2.16		Crippen Method
logp	2.239		Crippen Method
mvol	127.810	ml/mol	McGowan Method
pc	5183.17	kPa	Joback Method
tb	652.35	K	Joback Method
tc	905.23	K	Joback Method
tf	516.27	K	Joback Method
vc	0.384	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	352.76	J/molxK	652.35	Joback Method
cpg	402.60	J/molxK	863.08	Joback Method
cpg	393.19	J/molxK	820.94	Joback Method
cpg	383.78	J/molxK	778.79	Joback Method
cpg	374.09	J/molxK	736.64	Joback Method
cpg	363.84	J/molxK	694.50	Joback Method
cpg	412.28	J/molxK	905.23	Joback Method
dvisc	0.0000292	Paxs	652.35	Joback Method
dvisc	0.0000385	Paxs	629.67	Joback Method

dvisc	0.0000519	Paxs	606.99	Joback Method
dvisc	0.0000715	Paxs	584.31	Joback Method
dvisc	0.0001011	Paxs	561.63	Joback Method
dvisc	0.0001471	Paxs	538.95	Joback Method
dvisc	0.0002214	Paxs	516.27	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C3090479&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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
mccvol:	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/79-104-9/5-8-Endomethylene-5-8-dihydro-1-4-naphthohydroquinone.pdf>

Generated by Cheméo on 2024-05-04 05:11:31.237495393 +0000 UTC m=+17088740.158072705.

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