

Tricyclo[3.3.1.1^{3,7}]decane, tricyclo[3.3.1.1^{3,7}]decane

Other names:	[2,2']Biadamantanylidene 2,2'-Biadamantanyliden
Inchi:	InChI=1S/C20H28/c1-11-3-15-5-12(1)6-16(4-11)19(15)20-17-7-13-2-14(9-17)10-18(20)8
InchiKey:	KRDXURLKZNAXAQ-VXPUYCOJSA-N
Formula:	C20H28
SMILES:	C1C2CC3CC1CC(C2)C3=C1C2CC3CC(C2)CC1C3
Mol. weight [g/mol]:	268.44
CAS:	30541-56-1

Physical Properties

Property code	Value	Unit	Source
gf	453.10	kJ/mol	Joback Method
hf	-37.49	kJ/mol	Joback Method
hfus	34.75	kJ/mol	Joback Method
hvap	60.93	kJ/mol	Joback Method
ie	7.84	eV	NIST Webbook
log10ws	-5.48		Crippen Method
logp	5.195		Crippen Method
mcvol	223.200	ml/mol	McGowan Method
pc	1783.35	kPa	Joback Method
tb	705.76	K	Joback Method
tc	939.61	K	Joback Method
tf	433.08	K	Joback Method
vc	0.866	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	748.10	J/mol×K	705.76	Joback Method
cpg	860.77	J/mol×K	900.64	Joback Method
cpg	840.69	J/mol×K	861.66	Joback Method
cpg	819.62	J/mol×K	822.69	Joback Method
cpg	797.32	J/mol×K	783.71	Joback Method
cpg	773.55	J/mol×K	744.74	Joback Method

cpg	880.09	J/mol×K	939.61	Joback Method
dvisc	0.0329500	Paxs	705.76	Joback Method
dvisc	0.0294018	Paxs	660.31	Joback Method
dvisc	0.0257975	Paxs	614.87	Joback Method
dvisc	0.0221675	Paxs	569.42	Joback Method
dvisc	0.0185536	Paxs	523.97	Joback Method
dvisc	0.0150128	Paxs	478.53	Joback Method
dvisc	0.0116196	Paxs	433.08	Joback Method

Sources

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=C30541561&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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
ie:	Ionization energy
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

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