

Hexafluoro-2,2-bis-(3,4-dimethylphenyl)-propane

Other names:	Benzene, 1,1'-[2,2,2-trifluoro-1-(trifluoromethyl)ethylidene]bis[3,4-dimethyl-2,2-Bis-(3,4-dimethylphenyl)-hexafluoropropane 1,1'-[2,2,2-trifluoro-1-(trifluoromethyl)ethylidene]bis[3,4-dimethylbenzene]
Inchi:	InChI=1S/C19H18F6/c1-11-5-7-15(9-13(11)3)17(18(20,21)22,19(23,24)25)16-8-6-12(2)1
InchiKey:	GLFKFHJEFMLTOB-UHFFFAOYSA-N
Formula:	C19H18F6
SMILES:	<chem>Cc1ccc(C(c2ccc(C)c(C)c2)(C(F)(F)F)C(F)(F)F)cc1C</chem>
Mol. weight [g/mol]:	360.34
CAS:	65294-20-4

Physical Properties

Property code	Value	Unit	Source
gf	-864.94	kJ/mol	Joback Method
hf	-1211.22	kJ/mol	Joback Method
hfus	27.73	kJ/mol	Joback Method
hvap	56.30	kJ/mol	Joback Method
log10ws	-7.31		Crippen Method
logp	6.331		Crippen Method
mcvol	241.670	ml/mol	McGowan Method
pc	1420.78	kPa	Joback Method
tb	693.33	K	Joback Method
tc	894.59	K	Joback Method
tf	417.61	K	Joback Method
vc	0.959	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	685.56	J/molxK	693.33	Joback Method
cpg	701.54	J/molxK	726.87	Joback Method
cpg	716.37	J/molxK	760.42	Joback Method
cpg	730.15	J/molxK	793.96	Joback Method
cpg	742.96	J/molxK	827.50	Joback Method
cpg	754.90	J/molxK	861.05	Joback Method

cpg

766.07

J/mol×K

894.59

Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	388.00 ± 5.00	K	0.30	NIST Webbook

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=C65294204&Units=SI

Legend

cpg:	Ideal gas heat capacity
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
tbrp:	Boiling point at reduced pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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