

2,2',5,5'-Tetrabromobiphenyl

Inchi:	InChI=1S/C12H6Br4/c13-7-1-3-11(15)9(5-7)10-6-8(14)2-4-12(10)16/h1-6H
InchiKey:	XEFMFJLRXHQLEM-UHFFFAOYSA-N
Formula:	C12H6Br4
SMILES:	BrC1ccc(Br)c(-c2cc(Br)ccc2Br)c1
Mol. weight [g/mol]:	469.79
CAS:	59080-37-4

Physical Properties

Property code	Value	Unit	Source
gf	293.74	kJ/mol	Joback Method
hf	241.49	kJ/mol	Joback Method
hfus	34.50	kJ/mol	Joback Method
hvap	75.25	kJ/mol	Joback Method
log10ws	-8.06		Aqueous Solubility Prediction Method
logp	6.404		Crippen Method
mcvol	202.420	ml/mol	McGowan Method
pc	4492.23	kPa	Joback Method
tb	811.88	K	Joback Method
tc	1111.87	K	Joback Method
tf	567.12	K	Joback Method
vc	0.740	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	395.65	J/molxK	811.88	Joback Method
cpg	404.43	J/molxK	861.88	Joback Method
cpg	412.49	J/molxK	911.88	Joback Method
cpg	420.02	J/molxK	961.87	Joback Method
cpg	427.16	J/molxK	1011.87	Joback Method
cpg	434.09	J/molxK	1061.87	Joback Method
cpg	440.97	J/molxK	1111.87	Joback Method
dvisc	0.0004823	Paxs	567.12	Joback Method

dvisc	0.0003613	Paxs	607.91	Joback Method
dvisc	0.0002807	Paxs	648.71	Joback Method
dvisc	0.0002246	Paxs	689.50	Joback Method
dvisc	0.0001843	Paxs	730.29	Joback Method
dvisc	0.0001544	Paxs	771.09	Joback Method
dvisc	0.0001317	Paxs	811.88	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
Aqueous Solubility Prediction Method:	http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C59080374&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

Legend

cp_g:	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:	Log ₁₀ of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
m_cvol:	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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