1,3-Bis(3-aminophenoxy)benzene

Other names: Benzenamine, 3,3'-[1,3-phenylenebis(oxy)]bis-
3,3'-[m-phenylenebis(oxy)]dianiline
Inchi: InChI=1S/C18H16N2O2/c19-13-4-1-6-15(10-13)21-17-8-3-9-18(12-17)22-16-7-2-5-14(20
InchiKey: DKKYOQYISDAQER-UHFFFAOYSA-N
Formula: C18H16N2O2
SMILES: Nc1cccc(Oc2cccc(Oc3cccc(N)c3)c2)c1
Mol. weight [g/mol]: 292.33
CAS: 10526-07-5

Physical Properties

Property code	Value	Unit	Source
gf	331.92	kJ/mol	Joback Method
hf	63.47	kJ/mol	Joback Method
hfus	36.10	kJ/mol	Joback Method
hvap	90.58	kJ/mol	Joback Method
log10ws	-4.01		Crippen Method
logp	4.436		Crippen Method
mcvol	224.900	ml/mol	McGowan Method
pc	2718.33	kPa	Joback Method
tb	896.12	K	Joback Method
tc	1163.63	K	Joback Method
tf	620.42	K	Joback Method
vc	0.814	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	665.10	J/mol×K	896.12	Joback Method
cpg	677.74	J/mol×K	940.71	Joback Method
cpg	688.86	J/mol×K	985.29	Joback Method
cpg	698.53	J/mol×K	1029.88	Joback Method
cpg	706.80	J/mol×K	1074.46	Joback Method
cpg	713.75	J/mol×K	1119.05	Joback Method
cpg	719.42	J/mol×K	1163.63	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C10526075&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume