

Alpha,alpha,alpha,2,5-pentachloroacetanilide

Inchi:	InChI=1S/C8H4Cl5NO/c9-4-1-2-5(10)6(3-4)14-7(15)8(11,12)13/h1-3H,(H,14,15)
InchiKey:	WKBFRYPEBOWXGM-UHFFFAOYSA-N
Formula:	C8H4Cl5NO
SMILES:	O=C(Nc1cc(Cl)ccc1Cl)C(Cl)(Cl)Cl
Mol. weight [g/mol]:	307.39
CAS:	33715-64-9

Physical Properties

Property code	Value	Unit	Source
gf	13.29	kJ/mol	Joback Method
hf	-141.42	kJ/mol	Joback Method
hfus	30.01	kJ/mol	Joback Method
hvap	70.81	kJ/mol	Joback Method
log10ws	-4.61		Crippen Method
logp	4.302		Crippen Method
mcvol	172.570	ml/mol	McGowan Method
pc	3173.97	kPa	Joback Method
tb	707.04	K	Joback Method
tc	963.68	K	Joback Method
tf	485.99	K	Joback Method
vc	0.650	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	340.50	J/molxK	707.04	Joback Method
cpg	347.58	J/molxK	749.81	Joback Method
cpg	353.89	J/molxK	792.59	Joback Method
cpg	359.50	J/molxK	835.36	Joback Method
cpg	364.51	J/molxK	878.13	Joback Method
cpg	368.99	J/molxK	920.90	Joback Method
cpg	373.03	J/molxK	963.68	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
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=C33715649&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
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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