

Mexiletine

Other names:	2-Propanamine, 1-(2,6-dimethylphenoxy)- Ethylamine, 1-methyl-2-(2,6-xilyloxy)- 1-(2,6-Dimethylphenoxy)-2-aminopropane 1-(2',6'-Dimethylphenoxy)-2-aminopropane 1-Methyl-2-(2,6-xilyloxy)ethylamine 1-(2,6-Dimethylphenoxy)-2-propanamine Mexiletina 2-Amino-1-(2,6-dimethylphenoxy)propane 5370-01-4 Mexilitine
Inchi:	InChI=1S/C11H17NO/c1-8-5-4-6-9(2)11(8)13-7-10(3)12/h4-6,10H,7,12H2,1-3H3
InchiKey:	VLPIATFUUWWMKC-UHFFFAOYSA-N
Formula:	C11H17NO
SMILES:	<chem>Cc1cccc(C)c1OCC(C)N</chem>
Mol. weight [g/mol]:	179.26
CAS:	31828-71-4

Physical Properties

Property code	Value	Unit	Source
gf	93.90	kJ/mol	Joback Method
hf	-160.49	kJ/mol	Joback Method
hfus	20.37	kJ/mol	Joback Method
hvap	56.34	kJ/mol	Joback Method
log10ws	-2.93		Crippen Method
logp	2.029		Crippen Method
mcvol	157.940	ml/mol	McGowan Method
pc	2698.60	kPa	Joback Method
rinpol	1408.00		NIST Webbook
tb	582.23	K	Joback Method
tc	798.66	K	Joback Method
tf	355.68	K	Joback Method
vc	0.585	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	388.03	J/mol×K	582.23	Joback Method
cpg	403.25	J/mol×K	618.30	Joback Method
cpg	417.65	J/mol×K	654.37	Joback Method
cpg	431.25	J/mol×K	690.44	Joback Method
cpg	444.07	J/mol×K	726.51	Joback Method
cpg	456.12	J/mol×K	762.58	Joback Method
cpg	467.41	J/mol×K	798.66	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C31828714&Units=SI
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

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
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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