

SPIROXAMINE-1

Other names:	SPIROXAMINE-1 SPIROXAMINE-2 Spiroxamin 1,4-Dioxaspiro[4.5]decane-2-methanamine, 8-(1,1-dimethylethyl)-N-ethyl-N-propyl- 8-(1,1-dimethylethyl)-N-ethyl-N-propyl-1,4-dioxaspiro[4.5]decane-2-methanamine Spiroxamine, isomer 1
Inchi:	InChI=1S/C18H35NO2/c1-6-12-19(7-2)13-16-14-20-18(21-16)10-8-15(9-11-18)17(3,4)5/1
InchiKey:	PUYXTUJWRLOUCW-UHFFFAOYSA-N
Formula:	C18H35NO2
SMILES:	CCCN(CC)CC1COC2(CCC(C(C)(C)C)CC2)O1
Mol. weight [g/mol]:	297.48

Physical Properties

Property code	Value	Unit	Source
hfus	36.58	kJ/mol	Joback Method
logp	4.066		Crippen Method
mcvol	264.480	ml/mol	McGowan Method
rinpol	1896.00		NIST Webbook
rinpol	1896.00		NIST Webbook
rinpol	1896.00		NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	830.03	J/mol×K	700.48	Joback Method
cpg	854.02	J/mol×K	735.20	Joback Method
cpg	876.74	J/mol×K	769.93	Joback Method
cpg	898.36	J/mol×K	804.65	Joback Method
cpg	919.03	J/mol×K	839.38	Joback Method
cpg	938.93	J/mol×K	874.10	Joback Method
cpg	958.21	J/mol×K	908.83	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C118134308&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	

Legend

cpg:	Ideal gas heat capacity
hfus:	Enthalpy of fusion at standard conditions
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices

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