Benzamide, 2-fluoro-N-tetradecyl-

Inchi: InChI=1S/C21H34FNO/c1-2-3-4-5-6-7-8-9-10-11-12-15-18-23-21(24)19-16-13-14-17-20(
InchiKey: VYZGAKXLSSMCLT-UHFFFAOYSA-N
Formula: C21H34FNO
SMILES: CCCCCCCCCCCCCCN=C(O)c1ccccc1F
Mol. weight [g/mol]: 335.50

Physical Properties

Property code	Value	Unit	Source
hf	-527.62	kJ/mol	Joback Method
hvap	84.53	kJ/mol	Joback Method
log10ws	-7.13		Crippen Method
logp	6.831		Crippen Method
mcvol	296.310	ml/mol	McGowan Method
pc	1132.91	kPa	Joback Method
rinpol	2647.00		NIST Webbook
rinpol	2647.00		NIST Webbook
tb	879.55	K	Joback Method
tc	1078.81	K	Joback Method

Sources

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
McGowan Method: http://link.springer.com/article/10.1007/BF02311772
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U407143&Units=SI

Legend

hf: Enthalpy of formation 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

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