

4-Fluoro-3-trifluoromethylbenzoic acid, heptadecyl ester

Inchi:	InChI=1S/C25H38F4O2/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-16-19-31-24(30)21-17-18
InchiKey:	NYJXVRCADIFCSE-UHFFFAOYSA-N
Formula:	C25H38F4O2
SMILES:	CCCCCCCCCCCCCCCCOC(=O)c1ccc(F)c(C(F)(F)F)c1
Mol. weight [g/mol]:	446.56

Physical Properties

Property code	Value	Unit	Source
gf	-757.55	kJ/mol	Joback Method
hf	-1383.73	kJ/mol	Joback Method
hfus	61.46	kJ/mol	Joback Method
hvap	79.44	kJ/mol	Joback Method
log10ws	-9.85		Crippen Method
logp	8.873		Crippen Method
mvol	353.870	ml/mol	McGowan Method
pc	849.00	kPa	Joback Method
rinpol	2620.00		NIST Webbook
rinpol	2620.00		NIST Webbook
tb	878.18	K	Joback Method
tc	1075.21	K	Joback Method
tf	499.91	K	Joback Method
vc	1.413	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1158.56	J/molxK	878.18	Joback Method
cpg	1177.13	J/molxK	911.02	Joback Method
cpg	1194.54	J/molxK	943.86	Joback Method
cpg	1210.84	J/molxK	976.70	Joback Method
cpg	1226.09	J/molxK	1009.53	Joback Method
cpg	1240.36	J/molxK	1042.37	Joback Method
cpg	1253.72	J/molxK	1075.21	Joback Method

Sources

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=U338929&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I

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

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