

4-Fluoro-2-trifluoromethylbenzoic acid, neopentyl ester

Inchi:	InChI=1S/C13H14F4O2/c1-12(2,3)7-19-11(18)9-5-4-8(14)6-10(9)13(15,16)17/h4-6H,7H2
InchiKey:	FYMAZHRUKGZZEQ-UHFFFAOYSA-N
Formula:	C13H14F4O2
SMILES:	CC(C)(C)COC(=O)c1ccc(F)cc1C(F)(F)F
Mol. weight [g/mol]:	278.24

Physical Properties

Property code	Value	Unit	Source
gf	-855.75	kJ/mol	Joback Method
hf	-1144.80	kJ/mol	Joback Method
hfus	22.97	kJ/mol	Joback Method
hvap	51.43	kJ/mol	Joback Method
log10ws	-4.58		Crippen Method
logp	4.047		Crippen Method
mcvol	184.790	ml/mol	McGowan Method
pc	1956.14	kPa	Joback Method
rinqol	1288.00		NIST Webbook
tb	600.39	K	Joback Method
tc	791.41	K	Joback Method
tf	367.09	K	Joback Method
vc	0.730	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	481.60	J/molxK	600.39	Joback Method
cpg	495.65	J/molxK	632.23	Joback Method
cpg	508.79	J/molxK	664.06	Joback Method
cpg	521.09	J/molxK	695.90	Joback Method
cpg	532.59	J/molxK	727.74	Joback Method
cpg	543.32	J/molxK	759.58	Joback Method
cpg	553.33	J/molxK	791.41	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=U357657&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
hvac:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mccvol:	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

Latest version available from:

<https://www.chemeo.com/cid/52-273-1/4-Fluoro-2-trifluoromethylbenzoic-acid-neopentyl-ester.pdf>

Generated by Cheméo on 2024-04-26 15:57:35.637422 +0000 UTC m=+16436304.557999312.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.