

6-Fluoro-2-trifluoromethylbenzoic acid, 2,2,3,3,4,4,4-heptafluorobutyl ester

Other names:	6-Fluoro-2-trifluorobenzoic acid, 2,2,3,3,4,4,4-heptafluorobutyl ester
Inchi:	InChI=1S/C12H5F11O2/c13-6-3-1-2-5(10(16,17)18)7(6)8(24)25-4-9(14,15)11(19,20)12(2
InchiKey:	OQANPNONJGUGRF-UHFFFAOYSA-N
Formula:	C12H5F11O2
SMILES:	O=C(OCC(F)(F)C(F)(F)C(F)(F)F)c1c(F)cccc1C(F)(F)F
Mol. weight [g/mol]:	390.15

Physical Properties

Property code	Value	Unit	Source
gf	-2222.16	kJ/mol	Joback Method
hf	-2514.43	kJ/mol	Joback Method
hfus	27.11	kJ/mol	Joback Method
hvap	40.89	kJ/mol	Joback Method
log10ws	-5.70		Crippen Method
logp	4.834		Crippen Method
mcvol	183.090	ml/mol	McGowan Method
pc	1676.91	kPa	Joback Method
rinpol	1121.00		NIST Webbook
rinpol	1121.00		NIST Webbook
tb	565.94	K	Joback Method
tc	725.76	K	Joback Method
tf	364.79	K	Joback Method
vc	0.777	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	496.91	J/molxK	565.94	Joback Method
cpg	508.07	J/molxK	592.58	Joback Method
cpg	518.41	J/molxK	619.21	Joback Method
cpg	528.00	J/molxK	645.85	Joback Method
cpg	536.88	J/molxK	672.49	Joback Method
cpg	545.08	J/molxK	699.12	Joback Method
cpg	552.67	J/molxK	725.76	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=U343734&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
hvp:	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
rinp:	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/120-111-4/6-Fluoro-2-trifluoromethylbenzoic-acid-2-2-3-3-4-4-4-heptafluorobutyl-ester.p>

Generated by Cheméo on 2024-05-05 19:21:12.560958776 +0000 UTC m=+17226121.481536098.

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