

6-Fluoro-2-trifluoromethylbenzoic acid, 3-fluorophenyl ester

Other names: 6-Fluoro-2-trifluorobenzoic acid, 3-fluorophenyl ester

Inchi: InChI=1S/C14H7F5O2/c15-8-3-1-4-9(7-8)21-13(20)12-10(14(17,18)19)5-2-6-11(12)16/h1

InchiKey: ZURLLWOTCBBJQS-UHFFFAOYSA-N

Formula: C14H7F5O2

SMILES: O=C(Oc1cccc(F)c1)c1c(F)cccc1C(F)(F)F

Mol. weight [g/mol]: 302.20

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -942.20 | kJ/mol | Joback Method |
| hf | -1127.74 | kJ/mol | Joback Method |
| hfus | 29.70 | kJ/mol | Joback Method |
| hvap | 57.07 | kJ/mol | Joback Method |
| log10ws | -5.33 | | Crippen Method |
| logp | 4.203 | | Crippen Method |
| mcvol | 176.890 | ml/mol | McGowan Method |
| pc | 2256.81 | kPa | Joback Method |
| rinpol | 1589.00 | | NIST Webbook |
| rinpol | 1589.00 | | NIST Webbook |
| tb | 657.43 | K | Joback Method |
| tc | 864.10 | K | Joback Method |
| tf | 415.47 | K | Joback Method |
| vc | 0.707 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 456.51 | J/mol×K | 657.43 | Joback Method |
| cpg | 468.38 | J/mol×K | 691.87 | Joback Method |
| cpg | 479.35 | J/mol×K | 726.32 | Joback Method |
| cpg | 489.48 | J/mol×K | 760.76 | Joback Method |
| cpg | 498.79 | J/mol×K | 795.21 | Joback Method |
| cpg | 507.35 | J/mol×K | 829.65 | Joback Method |
| cpg | 515.17 | J/mol×K | 864.10 | Joback Method |

Sources

| | |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| 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=U343737&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.cheméo.com/doc/models/crippen_log10ws |

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 |
| h vap: | 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 |
| r in pol: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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