

5-Chlorovaleric acid, pentadecyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C20H39ClO2/c1-2-3-4-5-6-7-8-9-10-11-12-13-16-19-23-20(22)17-14-15-18-21 |
| InchiKey: | FGQMMDMNLZQATFL-UHFFFAOYSA-N |
| Formula: | C20H39ClO2 |
| SMILES: | CCCCCCCCCCCCCCCCOC(=O)CCCCCI |
| Mol. weight [g/mol]: | 346.98 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -128.33 | kJ/mol | Joback Method |
| hf | -716.67 | kJ/mol | Joback Method |
| hfus | 54.54 | kJ/mol | Joback Method |
| hvap | 73.66 | kJ/mol | Joback Method |
| log10ws | -7.21 | | Crippen Method |
| logp | 7.030 | | Crippen Method |
| mvol | 312.340 | ml/mol | McGowan Method |
| pc | 1021.38 | kPa | Joback Method |
| rinpol | 2465.90 | | NIST Webbook |
| rinpol | 2465.90 | | NIST Webbook |
| tb | 770.72 | K | Joback Method |
| tc | 948.17 | K | Joback Method |
| tf | 417.24 | K | Joback Method |
| vc | 1.228 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 944.23 | J/mol×K | 770.72 | Joback Method |
| cpg | 1029.30 | J/mol×K | 918.60 | Joback Method |
| cpg | 1014.07 | J/mol×K | 889.02 | Joback Method |
| cpg | 997.98 | J/mol×K | 859.45 | Joback Method |
| cpg | 980.99 | J/mol×K | 829.87 | Joback Method |
| cpg | 963.08 | J/mol×K | 800.30 | Joback Method |
| cpg | 1043.68 | J/mol×K | 948.17 | Joback Method |
| dvisc | 0.0000640 | Paxs | 770.72 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000858 | Paxs | 711.81 | Joback Method |
| dvisc | 0.0001213 | Paxs | 652.89 | Joback Method |
| dvisc | 0.0001837 | Paxs | 593.98 | Joback Method |
| dvisc | 0.0003049 | Paxs | 535.07 | Joback Method |
| dvisc | 0.0005737 | Paxs | 476.15 | Joback Method |
| dvisc | 0.0012902 | Paxs | 417.24 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U292316&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| 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 |
| rinpol: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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