

5-Bromovaleric acid, 2-pentadecyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C20H39BrO2/c1-3-4-5-6-7-8-9-10-11-12-13-16-19(2)23-20(22)17-14-15-18-21 |
| InchiKey: | HSOKXQMYYGQAEY-UHFFFAOYSA-N |
| Formula: | C20H39BrO2 |
| SMILES: | CCCCCCCCCCCCC(C)OC(=O)CCCCBr |
| Mol. weight [g/mol]: | 391.43 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -104.52 | kJ/mol | Joback Method |
| hf | -679.88 | kJ/mol | Joback Method |
| hfus | 52.10 | kJ/mol | Joback Method |
| hvap | 75.32 | kJ/mol | Joback Method |
| log10ws | -7.60 | | Crippen Method |
| logp | 7.184 | | Crippen Method |
| mvol | 317.600 | ml/mol | McGowan Method |
| pc | 1100.08 | kPa | Joback Method |
| rinpol | 2422.00 | | NIST Webbook |
| tb | 799.01 | K | Joback Method |
| tc | 983.68 | K | Joback Method |
| tf | 432.12 | K | Joback Method |
| vc | 1.236 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 963.67 | J/molxK | 799.01 | Joback Method |
| cpg | 1047.16 | J/molxK | 952.90 | Joback Method |
| cpg | 1032.29 | J/molxK | 922.13 | Joback Method |
| cpg | 1016.54 | J/molxK | 891.35 | Joback Method |
| cpg | 999.88 | J/molxK | 860.57 | Joback Method |
| cpg | 982.27 | J/molxK | 829.79 | Joback Method |
| cpg | 1061.18 | J/molxK | 983.68 | Joback Method |
| dvisc | 0.0000538 | Paxs | 799.01 | Joback Method |
| dvisc | 0.0000725 | Paxs | 737.86 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0001032 | Paxs | 676.71 | Joback Method |
| dvisc | 0.0001576 | Paxs | 615.57 | Joback Method |
| dvisc | 0.0002643 | Paxs | 554.42 | Joback Method |
| dvisc | 0.0005038 | Paxs | 493.27 | Joback Method |
| dvisc | 0.0011524 | Paxs | 432.12 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U299932&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 |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

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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