

2-Bromobenzoic acid, undec-2-enyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C18H25BrO2/c1-2-3-4-5-6-7-8-9-12-15-21-18(20)16-13-10-11-14-17(16)19/h9- |
| InchiKey: | JENAWSKCVBMMU-FMIVXFBMSA-N |
| Formula: | C18H25BrO2 |
| SMILES: | CCCCCCCC=CCOC(=O)c1ccccc1Br |
| Mol. weight [g/mol]: | 353.29 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 64.08 | kJ/mol | Joback Method |
| hf | -291.04 | kJ/mol | Joback Method |
| hfus | 44.30 | kJ/mol | Joback Method |
| hvap | 74.15 | kJ/mol | Joback Method |
| log10ws | -6.92 | | Crippen Method |
| logp | 5.913 | | Crippen Method |
| mcvol | 261.360 | ml/mol | McGowan Method |
| pc | 1648.43 | kPa | Joback Method |
| rinpol | 2327.00 | | NIST Webbook |
| rinpol | 2327.00 | | NIST Webbook |
| tb | 789.51 | K | Joback Method |
| tc | 998.63 | K | Joback Method |
| tf | 458.44 | K | Joback Method |
| vc | 1.002 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 731.35 | J/molxK | 789.51 | Joback Method |
| cpg | 800.58 | J/molxK | 963.78 | Joback Method |
| cpg | 788.44 | J/molxK | 928.92 | Joback Method |
| cpg | 775.51 | J/molxK | 894.07 | Joback Method |
| cpg | 761.72 | J/molxK | 859.22 | Joback Method |
| cpg | 747.02 | J/molxK | 824.36 | Joback Method |
| cpg | 811.97 | J/molxK | 998.63 | Joback Method |
| dvisc | 0.0000696 | Paxs | 789.51 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000892 | Paxs | 734.33 | Joback Method |
| dvisc | 0.0001192 | Paxs | 679.15 | Joback Method |
| dvisc | 0.0001676 | Paxs | 623.97 | Joback Method |
| dvisc | 0.0002517 | Paxs | 568.80 | Joback Method |
| dvisc | 0.0004127 | Paxs | 513.62 | Joback Method |
| dvisc | 0.0007619 | Paxs | 458.44 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U299399&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307I |
| 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 |
| 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 |

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