

3-Bromo-2,6-dichloroanisole

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|-----------------------------|--|
| Inchi: | InChI=1S/C7H5BrCl2O/c1-11-7-5(9)3-2-4(8)6(7)10/h2-3H,1H3 |
| InchiKey: | ANLVZYYGYBXMIX-UHFFFAOYSA-N |
| Formula: | C7H5BrCl2O |
| SMILES: | COc1c(Cl)ccc(Br)c1Cl |
| Mol. weight [g/mol]: | 255.92 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -22.96 | kJ/mol | Joback Method |
| hf | -123.06 | kJ/mol | Joback Method |
| hfus | 21.63 | kJ/mol | Joback Method |
| hvap | 53.05 | kJ/mol | Joback Method |
| log10ws | -4.12 | | Crippen Method |
| logp | 3.764 | | Crippen Method |
| mcvol | 133.580 | ml/mol | McGowan Method |
| pc | 3782.33 | kPa | Joback Method |
| rinpol | 1518.00 | | NIST Webbook |
| rinpol | 1518.00 | | NIST Webbook |
| tb | 564.62 | K | Joback Method |
| tc | 808.13 | K | Joback Method |
| tf | 374.50 | K | Joback Method |
| vc | 0.497 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 234.46 | J/molxK | 564.62 | Joback Method |
| cpg | 242.60 | J/molxK | 605.20 | Joback Method |
| cpg | 250.24 | J/molxK | 645.79 | Joback Method |
| cpg | 257.38 | J/molxK | 686.37 | Joback Method |
| cpg | 264.05 | J/molxK | 726.96 | Joback Method |
| cpg | 270.24 | J/molxK | 767.54 | Joback Method |
| cpg | 275.96 | J/molxK | 808.13 | Joback Method |
| dvisc | 0.0009926 | Paxs | 374.50 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0007114 | Paxs | 406.19 | Joback Method |
| dvisc | 0.0005351 | Paxs | 437.87 | Joback Method |
| dvisc | 0.0004183 | Paxs | 469.56 | Joback Method |
| dvisc | 0.0003373 | Paxs | 501.25 | Joback Method |
| dvisc | 0.0002790 | Paxs | 532.93 | Joback Method |
| dvisc | 0.0002358 | Paxs | 564.62 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R323601&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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