

3-Bromo-2-hydroxybenzoic acid

| | |
|-----------------------------|--|
| Inchi: | InChI=1S/C7H5BrO3/c8-5-3-1-2-4(6(5)9)7(10)11/h1-3,9H,(H,10,11) |
| InchiKey: | BHPSQWRVKOPSOQ-UHFFFAOYSA-N |
| Formula: | C7H5BrO3 |
| SMILES: | O=C(O)c1cccc(Br)c1O |
| Mol. weight [g/mol]: | 217.02 |
| CAS: | 3883-95-2 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -295.20 | kJ/mol | Joback Method |
| hf | -378.54 | kJ/mol | Joback Method |
| hfus | 24.29 | kJ/mol | Joback Method |
| hvap | 76.99 | kJ/mol | Joback Method |
| log10ws | -2.24 | | Crippen Method |
| logp | 1.853 | | Crippen Method |
| mcvol | 116.540 | ml/mol | McGowan Method |
| pc | 6932.88 | kPa | Joback Method |
| tb | 684.05 | K | Joback Method |
| tc | 916.63 | K | Joback Method |
| tf | 489.86 | K | Joback Method |
| vc | 0.372 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 255.56 | J/molxK | 684.05 | Joback Method |
| cpg | 282.69 | J/molxK | 877.87 | Joback Method |
| cpg | 277.76 | J/molxK | 839.11 | Joback Method |
| cpg | 272.67 | J/molxK | 800.34 | Joback Method |
| cpg | 267.33 | J/molxK | 761.58 | Joback Method |
| cpg | 261.65 | J/molxK | 722.81 | Joback Method |
| cpg | 287.53 | J/molxK | 916.63 | Joback Method |
| dvisc | 0.0000092 | Paxs | 684.05 | Joback Method |
| dvisc | 0.0000139 | Paxs | 651.69 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000220 | Paxs | 619.32 | Joback Method |
| dvisc | 0.0000366 | Paxs | 586.96 | Joback Method |
| dvisc | 0.0000647 | Paxs | 554.59 | Joback Method |
| dvisc | 0.0001225 | Paxs | 522.23 | Joback Method |
| dvisc | 0.0002524 | Paxs | 489.86 | 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=C3883952&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

Latest version available from:

<https://www.chemeo.com/cid/66-710-0/3-Bromo-2-hydroxybenzoic-acid.pdf>

Generated by Cheméo on 2024-04-26 05:33:02.836747979 +0000 UTC m=+16398831.757325294.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.