

# Bromoxynil

|                             |   |
|-----------------------------|---|
| <b>Other names:</b>         | 2,6-Dibromo-4-cyanophenol<br>3,5-Dibromo-4-hydroxybenzotrile<br>3,5-Dibromo-4-hydroxyphenylcyanide<br>4-Cyano-2,6-dibromophenol<br>4-Hydroxy-3,5-dibromobenzotrile<br>Benzotrile, 3,5-dibromo-4-hydroxy-<br>Brittox<br>Brominal<br>Brominex<br>Brominil<br>Bromotril<br>Broxynil<br>Bucril<br>Buctril<br>Buctril industrial<br>Butilchlorofos<br>Certrol B<br>Chipco buctril<br>Chipco crab-kleen<br>ENT 20852<br>Labuctril<br>Labuctril 25<br>Litarol M<br>M&B 10,064<br>M&B 10731<br>MB 10064<br>ME4 Brominal<br>Nu-lawn weeder<br>Oxytril M<br>Pardner<br>S 2132<br>Sabre<br>Toplan<br>Torch |
| <b>Inchi:</b>               | InChI=1S/C7H3Br2NO/c8-5-1-4(3-10)2-6(9)7(5)11/h1-2,11H  |
| <b>InchiKey:</b>            | UPMXNNIRAGDFEH-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C7H3Br2NO   |
| <b>SMILES:</b>              | N#Cc1cc(Br)c(O)c(Br)c1  |
| <b>Mol. weight [g/mol]:</b> | 276.91  |
| <b>CAS:</b>                 | 1689-84-5   |

# Physical Properties

| Property code | Value         | Unit                 | Source                               |
|---------------|---------------|----------------------|--------------------------------------|
| gf            | 108.41        | kJ/mol               | Joback Method                        |
| hf            | 66.01         | kJ/mol               | Joback Method                        |
| hfus          | 25.01         | kJ/mol               | Joback Method                        |
| hvap          | 71.14         | kJ/mol               | Joback Method                        |
| log10ws       | -3.33         |                      | Aqueous Solubility Prediction Method |
| log10ws       | -3.33         |                      | Estimated Solubility Method          |
| logp          | 2.789         |                      | Crippen Method                       |
| mvol          | 127.980       | ml/mol               | McGowan Method                       |
| pc            | 5678.82       | kPa                  | Joback Method                        |
| rinpol        | 1652.00       |                      | NIST Webbook                         |
| rinpol        | 1686.00       |                      | NIST Webbook                         |
| rinpol        | 1652.00       |                      | NIST Webbook                         |
| tb            | 711.22        | K                    | Joback Method                        |
| tc            | 985.74        | K                    | Joback Method                        |
| tf            | 464.20 ± 0.20 | K                    | NIST Webbook                         |
| tf            | 463.60 ± 0.20 | K                    | NIST Webbook                         |
| vc            | 0.435         | m <sup>3</sup> /kmol | Joback Method                        |

# Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 241.09 | J/mol×K | 711.22          | Joback Method |
| cpg           | 246.39 | J/mol×K | 756.97          | Joback Method |
| cpg           | 251.39 | J/mol×K | 802.73          | Joback Method |
| cpg           | 256.23 | J/mol×K | 848.48          | Joback Method |
| cpg           | 261.05 | J/mol×K | 894.23          | Joback Method |
| cpg           | 266.00 | J/mol×K | 939.98          | Joback Method |
| cpg           | 271.20 | J/mol×K | 985.74          | Joback Method |
| hfust         | 32.03  | kJ/mol  | 464.00          | NIST Webbook  |
| hfust         | 32.03  | kJ/mol  | 464.00          | NIST Webbook  |

# Sources

|  |   |
|--|---|
| <b>Aqueous Solubility Prediction Method:</b> | <a href="http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa">http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa</a>     |
| <b>Estimated Solubility Method:</b>          | <a href="http://pubs.acs.org/doi/suppl/10.1021/ci034243x/suppl_file/ci034243xsi20040112_053635.txt">http://pubs.acs.org/doi/suppl/10.1021/ci034243x/suppl_file/ci034243xsi20040112_053635.txt</a> |
| <b>McGowan Method:</b>                       | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>   |
| <b>NIST Webbook:</b>                         | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C1689845&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1689845&amp;Units=SI</a>   |
| <b>Crippen Method:</b>                       | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>   |
| <b>Joback Method:</b>                        | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>   |

# Legend

|                 |   |
|-----------------|---|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hfust:</b>   | Enthalpy of fusion at a given temperature       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>rinpol:</b>  | Non-polar retention indices                     |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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