

# Bromoacetic acid, decyl ester

|                      |                                                                      |
|----------------------|----------------------------------------------------------------------|
| Other names:         | Decyl bromoacetate                                                   |
| Inchi:               | InChI=1S/C12H23BrO2/c1-2-3-4-5-6-7-8-9-10-15-12(14)11-13/h2-11H2,1H3 |
| InchiKey:            | OMRFZIDOMSMRSP-UHFFFAOYSA-N                                          |
| Formula:             | C12H23BrO2                                                           |
| SMILES:              | CCCCCCCCCCCCOC(=O)CBr                                                |
| Mol. weight [g/mol]: | 279.21                                                               |
| CAS:                 | 5436-93-1                                                            |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -169.44 | kJ/mol  | Joback Method  |
| hf            | -509.48 | kJ/mol  | Joback Method  |
| hfus          | 34.91   | kJ/mol  | Joback Method  |
| hvap          | 57.90   | kJ/mol  | Joback Method  |
| log10ws       | -4.14   |         | Crippen Method |
| logp          | 4.065   |         | Crippen Method |
| mcvol         | 204.880 | ml/mol  | McGowan Method |
| pc            | 1966.56 | kPa     | Joback Method  |
| rinpol        | 1684.80 |         | NIST Webbook   |
| tb            | 616.41  | K       | Joback Method  |
| tc            | 798.29  | K       | Joback Method  |
| tf            | 356.96  | K       | Joback Method  |
| vc            | 0.793   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 512.44 | J/mol×K | 616.41          | Joback Method |
| cpg           | 527.35 | J/mol×K | 646.72          | Joback Method |
| cpg           | 541.57 | J/mol×K | 677.04          | Joback Method |
| cpg           | 555.13 | J/mol×K | 707.35          | Joback Method |
| cpg           | 568.03 | J/mol×K | 737.67          | Joback Method |
| cpg           | 580.30 | J/mol×K | 767.98          | Joback Method |
| cpg           | 591.95 | J/mol×K | 798.29          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0020531 | Pa×s | 356.96 | Joback Method |
| dvisc | 0.0010894 | Pa×s | 400.20 | Joback Method |
| dvisc | 0.0006541 | Pa×s | 443.44 | Joback Method |
| dvisc | 0.0004300 | Pa×s | 486.68 | Joback Method |
| dvisc | 0.0003027 | Pa×s | 529.93 | Joback Method |
| dvisc | 0.0002247 | Pa×s | 573.17 | Joback Method |
| dvisc | 0.0001739 | Pa×s | 616.41 | Joback Method |

## Sources

|                        |                                                                                                                                             |
|------------------------|---------------------------------------------------------------------------------------------------------------------------------------------|
| <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=C5436931&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C5436931&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                   |
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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>dvisc:</b>   | Dynamic viscosity                               |
| <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>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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