

# 4-Bromobutyric acid, 3-ethylphenyl ester

|                      |                                                                                   |
|----------------------|-----------------------------------------------------------------------------------|
| Inchi:               | InChI=1S/C12H15BrO2/c1-2-10-5-3-6-11(9-10)15-12(14)7-4-8-13/h3,5-6,9H,2,4,7-8H2,1 |
| InchiKey:            | JNAWIEDHUGQKPZ-UHFFFAOYSA-N                                                       |
| Formula:             | C12H15BrO2                                                                        |
| SMILES:              | CCc1cccc(OC(=O)CCCBrc1                                                            |
| Mol. weight [g/mol]: | 271.15                                                                            |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -66.66  | kJ/mol  | Joback Method  |
| hf            | -284.42 | kJ/mol  | Joback Method  |
| hfus          | 28.56   | kJ/mol  | Joback Method  |
| hvap          | 60.84   | kJ/mol  | Joback Method  |
| log10ws       | -3.86   |         | Crippen Method |
| logp          | 3.329   |         | Crippen Method |
| mcvol         | 181.120 | ml/mol  | McGowan Method |
| pc            | 2659.77 | kPa     | Joback Method  |
| rinpol        | 1846.00 |         | NIST Webbook   |
| tb            | 648.07  | K       | Joback Method  |
| tc            | 867.01  | K       | Joback Method  |
| tf            | 395.90  | K       | Joback Method  |
| vc            | 0.685   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 433.91    | J/mol×K | 648.07          | Joback Method |
| cpg           | 494.20    | J/mol×K | 830.52          | Joback Method |
| cpg           | 483.75    | J/mol×K | 794.03          | Joback Method |
| cpg           | 472.52    | J/mol×K | 757.54          | Joback Method |
| cpg           | 460.50    | J/mol×K | 721.05          | Joback Method |
| cpg           | 447.63    | J/mol×K | 684.56          | Joback Method |
| cpg           | 503.90    | J/mol×K | 867.01          | Joback Method |
| dvisc         | 0.0001717 | Pa×s    | 648.07          | Joback Method |
| dvisc         | 0.0002144 | Pa×s    | 606.04          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002768 | Pa×s | 564.01 | Joback Method |
| dvisc | 0.0003722 | Pa×s | 521.99 | Joback Method |
| dvisc | 0.0005273 | Pa×s | 479.96 | Joback Method |
| dvisc | 0.0007985 | Pa×s | 437.93 | Joback Method |
| dvisc | 0.0013207 | Pa×s | 395.90 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U354689&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U354689&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>                                     |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</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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