

# Benzyl 2-hydroxy-6-methoxybenzoate

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C15H14O4/c1-18-13-9-5-8-12(16)14(13)15(17)19-10-11-6-3-2-4-7-11/h2-9,16H

CGNJMCLXIMWKDY-UHFFFAOYSA-N

C15H14O4

COc1cccc(O)c1C(=O)OCc1ccccc1

258.27

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -202.93 | kJ/mol  | Joback Method  |
| hf            | -445.67 | kJ/mol  | Joback Method  |
| hfus          | 32.06   | kJ/mol  | Joback Method  |
| hvap          | 78.78   | kJ/mol  | Joback Method  |
| log10ws       | -3.49   |         | Crippen Method |
| logp          | 2.758   |         | Crippen Method |
| mcvol         | 193.870 | ml/mol  | McGowan Method |
| pc            | 2976.29 | kPa     | Joback Method  |
| rinpol        | 2046.00 |         | NIST Webbook   |
| tb            | 780.27  | K       | Joback Method  |
| tc            | 1022.10 | K       | Joback Method  |
| tf            | 530.28  | K       | Joback Method  |
| vc            | 0.667   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 540.38    | J/mol×K | 780.27          | Joback Method |
| cpg           | 553.49    | J/mol×K | 820.57          | Joback Method |
| cpg           | 565.67    | J/mol×K | 860.88          | Joback Method |
| cpg           | 576.97    | J/mol×K | 901.18          | Joback Method |
| cpg           | 587.49    | J/mol×K | 941.49          | Joback Method |
| cpg           | 597.31    | J/mol×K | 981.79          | Joback Method |
| cpg           | 606.51    | J/mol×K | 1022.10         | Joback Method |
| dvisc         | 0.0001200 | Pa×s    | 530.28          | Joback Method |
| dvisc         | 0.0000625 | Pa×s    | 571.95          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000356 | Pa×s | 613.61 | Joback Method |
| dvisc | 0.0000218 | Pa×s | 655.28 | Joback Method |
| dvisc | 0.0000141 | Pa×s | 696.94 | Joback Method |
| dvisc | 0.0000096 | Pa×s | 738.61 | Joback Method |
| dvisc | 0.0000068 | Pa×s | 780.27 | 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=R283588&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R283588&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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