

# Benzoic acid, 2,3-dihydroxy-, methyl ester

|                      |                                                                    |
|----------------------|--------------------------------------------------------------------|
| Other names:         | o-Pyrocatechuic acid, methyl ester<br>Methyl 2,3-dihydroxybenzoate |
| Inchi:               | InChI=1S/C8H8O4/c1-12-8(11)5-3-2-4-6(9)7(5)10/h2-4,9-10H,1H3       |
| InchiKey:            | DOAJWTSNTNAEIY-UHFFFAOYSA-N                                        |
| Formula:             | C8H8O4                                                             |
| SMILES:              | COC(=O)c1cccc(O)c1O                                                |
| Mol. weight [g/mol]: | 168.15                                                             |
| CAS:                 | 2411-83-8                                                          |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -414.27 | kJ/mol  | Joback Method  |
| hf            | -571.34 | kJ/mol  | Joback Method  |
| hfus          | 24.87   | kJ/mol  | Joback Method  |
| hvap          | 70.86   | kJ/mol  | Joback Method  |
| log10ws       | -0.81   |         | Crippen Method |
| logp          | 0.884   |         | Crippen Method |
| mcvol         | 119.000 | ml/mol  | McGowan Method |
| pc            | 5818.28 | kPa     | Joback Method  |
| tb            | 646.65  | K       | Joback Method  |
| tc            | 888.33  | K       | Joback Method  |
| tf            | 501.94  | K       | Joback Method  |
| vc            | 0.332   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 299.69 | J/mol×K | 646.65          | Joback Method |
| cpg           | 308.47 | J/mol×K | 686.93          | Joback Method |
| cpg           | 316.66 | J/mol×K | 727.21          | Joback Method |
| cpg           | 324.37 | J/mol×K | 767.49          | Joback Method |
| cpg           | 331.74 | J/mol×K | 807.77          | Joback Method |
| cpg           | 338.88 | J/mol×K | 848.05          | Joback Method |
| cpg           | 345.91 | J/mol×K | 888.33          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000678 | Pa×s | 501.94 | Joback Method |
| dvisc | 0.0000375 | Pa×s | 526.06 | Joback Method |
| dvisc | 0.0000218 | Pa×s | 550.18 | Joback Method |
| dvisc | 0.0000133 | Pa×s | 574.29 | Joback Method |
| dvisc | 0.0000084 | Pa×s | 598.41 | Joback Method |
| dvisc | 0.0000055 | Pa×s | 622.53 | Joback Method |
| dvisc | 0.0000038 | Pa×s | 646.65 | Joback Method |

## Sources

|                        |                                                                                                                                             |
|------------------------|---------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C2411838&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C2411838&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>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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