

# Phenol, 3-methyl-2-iodo

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C7H7IO/c1-5-3-2-4-6(9)7(5)8/h2-4,9H,1H3 |
| <b>InchiKey:</b>            | SQQLELZQLWQUEP-UHFFFAOYSA-N                      |
| <b>Formula:</b>             | C7H7IO   |
| <b>SMILES:</b>              | Cc1cccc(O)c1I                                    |
| <b>Mol. weight [g/mol]:</b> | 234.03   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 14.34   | kJ/mol               | Joback Method  |
| hf            | -63.19  | kJ/mol               | Joback Method  |
| hfus          | 17.73   | kJ/mol               | Joback Method  |
| hvap          | 56.50   | kJ/mol               | Joback Method  |
| log10ws       | -2.64   |                      | Crippen Method |
| logp          | 2.305   |                      | Crippen Method |
| mcvol         | 117.420 | ml/mol               | McGowan Method |
| pc            | 4775.98 | kPa                  | Joback Method  |
| rinpol        | 1262.00 |                      | NIST Webbook   |
| rinpol        | 1262.00 |                      | NIST Webbook   |
| tb            | 564.98  | K                    | Joback Method  |
| tc            | 831.06  | K                    | Joback Method  |
| tf            | 377.37  | K                    | Joback Method  |
| vc            | 0.373   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 222.73    | J/molxK | 564.98          | Joback Method |
| cpg           | 231.80    | J/molxK | 609.33          | Joback Method |
| cpg           | 240.04    | J/molxK | 653.67          | Joback Method |
| cpg           | 247.59    | J/molxK | 698.02          | Joback Method |
| cpg           | 254.56    | J/molxK | 742.37          | Joback Method |
| cpg           | 261.10    | J/molxK | 786.71          | Joback Method |
| cpg           | 267.33    | J/molxK | 831.06          | Joback Method |
| dvisc         | 0.0019284 | Paxs    | 377.37          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0008888 | Paxs | 408.64 | Joback Method |
| dvisc | 0.0004574 | Paxs | 439.91 | Joback Method |
| dvisc | 0.0002570 | Paxs | 471.18 | Joback Method |
| dvisc | 0.0001552 | Paxs | 502.44 | Joback Method |
| dvisc | 0.0000994 | Paxs | 533.71 | Joback Method |
| dvisc | 0.0000669 | Paxs | 564.98 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R632554&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R632554&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>                         |

## 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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