

# 4-Butylbenzoic acid, tetradecyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C25H42O2/c1-3-5-7-8-9-10-11-12-13-14-15-16-22-27-25(26)24-20-18-23(19-2 |
| <b>InchiKey:</b>            | KTDIJHFLBJNYQD-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C25H42O2   |
| <b>SMILES:</b>              | CCCCCCCCCCCCCOC(=O)c1ccc(CCCC)cc1  |
| <b>Mol. weight [g/mol]:</b> | 374.60   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 28.48   | kJ/mol  | Joback Method  |
| hf            | -579.07 | kJ/mol  | Joback Method  |
| hfus          | 56.94   | kJ/mol  | Joback Method  |
| hvap          | 83.34   | kJ/mol  | Joback Method  |
| log10ws       | -8.80   |         | Crippen Method |
| logp          | 7.887   |         | Crippen Method |
| mcvol         | 346.790 | ml/mol  | McGowan Method |
| pc            | 943.26  | kPa     | Joback Method  |
| rinpol        | 2754.70 |         | NIST Webbook   |
| tb            | 879.35  | K       | Joback Method  |
| tc            | 1078.25 | K       | Joback Method  |
| tf            | 482.61  | K       | Joback Method  |
| vc            | 1.351   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1129.80   | J/molxK | 879.35          | Joback Method |
| cpg           | 1149.32   | J/molxK | 912.50          | Joback Method |
| cpg           | 1167.61   | J/molxK | 945.65          | Joback Method |
| cpg           | 1184.71   | J/molxK | 978.80          | Joback Method |
| cpg           | 1200.69   | J/molxK | 1011.95         | Joback Method |
| cpg           | 1215.57   | J/molxK | 1045.10         | Joback Method |
| cpg           | 1229.43   | J/molxK | 1078.25         | Joback Method |
| dvisc         | 0.0006292 | Paxs    | 482.61          | Joback Method |
| dvisc         | 0.0002964 | Paxs    | 548.73          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001641 | Paxs | 614.86 | Joback Method |
| dvisc | 0.0001020 | Paxs | 680.98 | Joback Method |
| dvisc | 0.0000689 | Paxs | 747.10 | Joback Method |
| dvisc | 0.0000496 | Paxs | 813.23 | Joback Method |
| dvisc | 0.0000376 | Paxs | 879.35 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U292327&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U292327&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</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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