

# 1-Naphthaleneacetic acid, heptyl ester

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C19H24O2/c1-2-3-4-5-8-14-21-19(20)15-17-12-9-11-16-10-6-7-13-18(16)17/h6 |
| <b>InchiKey:</b>            | AQOHGCUPDSXLGK-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C19H24O2  |
| <b>SMILES:</b>              | CCCCCCCOC(=O)Cc1cccc2ccccc12  |
| <b>Mol. weight [g/mol]:</b> | 284.39  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 84.61   | kJ/mol               | Joback Method  |
| hf            | -264.16 | kJ/mol               | Joback Method  |
| hfus          | 38.42   | kJ/mol               | Joback Method  |
| hvap          | 71.62   | kJ/mol               | Joback Method  |
| log10ws       | -5.87   |                      | Crippen Method |
| logp          | 4.896   |                      | Crippen Method |
| mcvol         | 242.790 | ml/mol               | McGowan Method |
| pc            | 1686.56 | kPa                  | Joback Method  |
| rinsol        | 2277.00 |                      | NIST Webbook   |
| tb            | 761.05  | K                    | Joback Method  |
| tc            | 970.67  | K                    | Joback Method  |
| tf            | 447.69  | K                    | Joback Method  |
| vc            | 0.938   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 700.90    | J/molxK | 761.05          | Joback Method |
| cpg           | 717.36    | J/molxK | 795.99          | Joback Method |
| cpg           | 732.80    | J/molxK | 830.92          | Joback Method |
| cpg           | 747.25    | J/molxK | 865.86          | Joback Method |
| cpg           | 760.78    | J/molxK | 900.80          | Joback Method |
| cpg           | 773.46    | J/molxK | 935.73          | Joback Method |
| cpg           | 785.33    | J/molxK | 970.67          | Joback Method |
| dvisc         | 0.0011478 | Paxs    | 447.69          | Joback Method |
| dvisc         | 0.0006946 | Paxs    | 499.92          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0004622 | Paxs | 552.14 | Joback Method |
| dvisc | 0.0003300 | Paxs | 604.37 | Joback Method |
| dvisc | 0.0002486 | Paxs | 656.60 | Joback Method |
| dvisc | 0.0001953 | Paxs | 708.82 | Joback Method |
| dvisc | 0.0001585 | Paxs | 761.05 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U394003&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U394003&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>                                 |

## 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                                 |

Latest version available from:

<https://www.chemeo.com/cid/89-497-3/1-Naphthaleneacetic-acid-heptyl-ester.pdf>

Generated by Cheméo on 2024-04-23 14:37:12.733975985 +0000 UTC m=+16172281.654553303.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.