

# Heptadecyl trifluoroacetate

|                      |                                                                                                                   |
|----------------------|-------------------------------------------------------------------------------------------------------------------|
| Other names:         | Heptadecyl 2,2,2-trifluoroacetate<br>1-Heptadecanol, trifluoroacetate<br>Trifluoroacetic acid, n-heptadecyl ester |
| Inchi:               | InChI=1S/C19H35F3O2/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-16-17-24-18(23)19(20,21)                                 |
| InchiKey:            | CDOYOXVOYRQINU-UHFFFAOYSA-N                                                                                       |
| Formula:             | C19H35F3O2                                                                                                        |
| SMILES:              | CCCCCCCCCCCCCCCCCOC(=O)C(F)(F)F                                                                                   |
| Mol. weight [g/mol]: | 352.48                                                                                                            |

## Physical Properties

| Property code | Value    | Unit    | Source         |
|---------------|----------|---------|----------------|
| gf            | -706.41  | kJ/mol  | Joback Method  |
| hf            | -1277.37 | kJ/mol  | Joback Method  |
| hfus          | 49.58    | kJ/mol  | Joback Method  |
| hvap          | 63.30    | kJ/mol  | Joback Method  |
| log10ws       | -7.30    |         | Crippen Method |
| logp          | 6.963    |         | Crippen Method |
| mcvol         | 291.320  | ml/mol  | McGowan Method |
| pc            | 1032.57  | kPa     | Joback Method  |
| rinpol        | 1910.20  |         | NIST Webbook   |
| rinpol        | 1907.00  |         | NIST Webbook   |
| tb            | 704.99   | K       | Joback Method  |
| tc            | 868.50   | K       | Joback Method  |
| tf            | 380.24   | K       | Joback Method  |
| vc            | 1.167    | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 873.58 | J/mol×K | 704.99          | Joback Method |
| cpg           | 891.77 | J/mol×K | 732.24          | Joback Method |
| cpg           | 909.12 | J/mol×K | 759.49          | Joback Method |
| cpg           | 925.64 | J/mol×K | 786.74          | Joback Method |
| cpg           | 941.36 | J/mol×K | 813.99          | Joback Method |

|     |        |         |        |               |
|-----|--------|---------|--------|---------------|
| cpg | 956.32 | J/mol×K | 841.24 | Joback Method |
| cpg | 970.55 | J/mol×K | 868.50 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U351870&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U351870&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>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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