

# Formic acid, hept-2-yl ester

|                      |                                                          |
|----------------------|----------------------------------------------------------|
| Inchi:               | InChI=1S/C8H16O2/c1-3-4-5-6-8(2)10-7-9/h7-8H,3-6H2,1-2H3 |
| InchiKey:            | SEIYVCWCDDKKIS-UHFFFAOYSA-N                              |
| Formula:             | C8H16O2                                                  |
| SMILES:              | CCCCC(C)OC=O                                             |
| Mol. weight [g/mol]: | 144.21                                                   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -190.48 | kJ/mol  | Joback Method  |
| hf            | -431.53 | kJ/mol  | Joback Method  |
| hfus          | 16.43   | kJ/mol  | Joback Method  |
| hvap          | 42.14   | kJ/mol  | Joback Method  |
| log10ws       | -2.15   |         | Crippen Method |
| logp          | 2.128   |         | Crippen Method |
| mcvol         | 131.020 | ml/mol  | McGowan Method |
| pc            | 2709.85 | kPa     | Joback Method  |
| rinpol        | 986.00  |         | NIST Webbook   |
| tb            | 453.08  | K       | Joback Method  |
| tc            | 627.37  | K       | Joback Method  |
| tf            | 229.15  | K       | Joback Method  |
| vc            | 0.512   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 280.99    | J/mol×K | 453.08          | Joback Method |
| cpg           | 337.83    | J/mol×K | 598.32          | Joback Method |
| cpg           | 327.32    | J/mol×K | 569.27          | Joback Method |
| cpg           | 316.38    | J/mol×K | 540.23          | Joback Method |
| cpg           | 305.02    | J/mol×K | 511.18          | Joback Method |
| cpg           | 293.22    | J/mol×K | 482.13          | Joback Method |
| cpg           | 347.93    | J/mol×K | 627.37          | Joback Method |
| dvisc         | 0.0002610 | Pa×s    | 453.08          | Joback Method |
| dvisc         | 0.0003456 | Pa×s    | 415.76          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0004836 | Pa×s | 378.44 | Joback Method |
| dvisc | 0.0007283 | Pa×s | 341.12 | Joback Method |
| dvisc | 0.0012128 | Pa×s | 303.79 | Joback Method |
| dvisc | 0.0023301 | Pa×s | 266.47 | Joback Method |
| dvisc | 0.0055375 | Pa×s | 229.15 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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=U368692&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U368692&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>                                     |

## 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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