

# 4-Fluoromethoxyacetophenone

|                      |                                                                                     |
|----------------------|-------------------------------------------------------------------------------------|
| Other names:         | 1-(4-Trifluoromethoxy-phenyl)-ethanone<br>1-[4-(trifluoromethoxy)phenyl]ethan-1-one |
| Inchi:               | InChI=1S/C9H7F3O2/c1-6(13)7-2-4-8(5-3-7)14-9(10,11)12/h2-5H,1H3                     |
| InchiKey:            | MOEXTBIPPMLEFX-UHFFFAOYSA-N                                                         |
| Formula:             | C9H7F3O2                                                                            |
| SMILES:              | CC(=O)c1ccc(OC(F)(F)F)cc1                                                           |
| Mol. weight [g/mol]: | 204.15                                                                              |

## Physical Properties

| Property code | Value   | Unit    | Source             |
|---------------|---------|---------|--------------------|
| af            | 0.5103  |         | Cheméo Relay (1.0) |
| gf            | -687.83 | kJ/mol  | Joback Method      |
| hf            | -945.87 | kJ/mol  | Cheméo Relay (1.0) |
| hfus          | 17.33   | kJ/mol  | Joback Method      |
| hvap          | 57.21   | kJ/mol  | Cheméo Relay (1.0) |
| ie            | 9.55    | eV      | Cheméo Relay (1.0) |
| log10ws       | -2.90   |         | Cheméo Relay (1.0) |
| logp          | 2.788   |         | Crippen Method     |
| mcvol         | 126.660 | ml/mol  | McGowan Method     |
| pc            | 2966.57 | kPa     | Joback Method      |
| tb            | 465.43  | K       | Cheméo Relay (1.0) |
| tc            | 637.65  | K       | Cheméo Relay (1.0) |
| tf            | 274.08  | K       | Cheméo Relay (1.0) |
| vc            | 0.466   | m3/kmol | Cheméo Relay (1.0) |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 284.63 | J/mol×K | 507.85          | Joback Method |
| cpg           | 295.99 | J/mol×K | 540.87          | Joback Method |
| cpg           | 306.64 | J/mol×K | 573.90          | Joback Method |
| cpg           | 316.61 | J/mol×K | 606.92          | Joback Method |
| cpg           | 325.92 | J/mol×K | 639.95          | Joback Method |
| cpg           | 334.60 | J/mol×K | 672.97          | Joback Method |

## Sources

|                           |                                                                                                                                               |
|---------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| McGowan Method:           | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                         |
| Crippen Method:           | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                     |
| Crippen Method:           | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                             |
| Cheméo Relay (1.0):       |                                                                                                                                               |
| Cheméo Relay (1.0, beta): |                                                                                                                                               |
| NIST Webbook:             | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C85013985&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C85013985&amp;Units=SI</a> |
| Joback Method:            | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                         |

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>af:</b>      | Acentric Factor                                 |
| <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>ie:</b>      | Ionization energy                               |
| <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>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/65-797-6/4-Fluoromethoxyacetophenone.pdf>

Generated by Cheméo on 2026-07-21 16:40:58.972216467 +0000 UTC m=+163154.814101850.

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