

# 2-Trimethylsilylthiophene

|                      |                                                                                    |
|----------------------|------------------------------------------------------------------------------------|
| Other names:         | 2-Thienyltrimethylsilane<br>2-Trimethylsilylthiophene<br>Trimethyl-2-thienylsilane |
| Inchi:               | InChI=1S/C7H12SSi/c1-9(2,3)7-5-4-6-8-7/h4-6H,1-3H3                                 |
| InchiKey:            | OANGLGSZPSFVDY-UHFFFAOYSA-N                                                        |
| Formula:             | C7H12SSi                                                                           |
| SMILES:              | C[Si](C)(C)c1cccs1                                                                 |
| Mol. weight [g/mol]: | 156.33                                                                             |

## Physical Properties

| Property code | Value   | Unit    | Source                   |
|---------------|---------|---------|--------------------------|
| hf            | -90.09  | kJ/mol  | Cheméo Relay (1.0)       |
| hvap          | 46.52   | kJ/mol  | Cheméo Relay (1.0)       |
| ie            | 8.10    | eV      | NIST Webbook             |
| ie            | 8.68    | eV      | NIST Webbook             |
| log10ws       | -3.22   |         | Cheméo Relay (1.0)       |
| logp          | 2.293   |         | Crippen Method           |
| pc            | 2696.38 | kPa     | Cheméo Relay (1.0, beta) |
| tb            | 440.47  | K       | Cheméo Relay (1.0)       |
| tc            | 640.39  | K       | Cheméo Relay (1.0)       |
| tf            | 213.27  | K       | Cheméo Relay (1.0)       |
| vc            | 0.479   | m3/kmol | Cheméo Relay (1.0)       |

## Sources

|                           |                                                                                                                                               |
|---------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| NIST Webbook:             | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C18245288&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C18245288&amp;Units=SI</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): |                                                                                                                                               |

# Legend

|                            |                                                 |
|----------------------------|-------------------------------------------------|
| <b>hf:</b>                 | Enthalpy of formation at standard conditions    |
| <b>h<sub>vap</sub>:</b>    | Enthalpy of vaporization at standard conditions |
| <b>ie:</b>                 | Ionization energy                               |
| <b>log<sub>10</sub>ws:</b> | Log10 of Water solubility in mol/l              |
| <b>log<sub>p</sub>:</b>    | Octanol/Water partition coefficient             |
| <b>p<sub>c</sub>:</b>      | Critical Pressure                               |
| <b>t<sub>b</sub>:</b>      | Normal Boiling Point Temperature                |
| <b>t<sub>c</sub>:</b>      | Critical Temperature                            |
| <b>t<sub>f</sub>:</b>      | Normal melting (fusion) point                   |
| <b>v<sub>c</sub>:</b>      | Critical Volume                                 |

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