

Sulfur monoxide

Inchi: InChI=1S/OS/c1-2
InchiKey: XTQHKBHJIVJGKJ-UHFFFAOYSA-N
Formula: OS
SMILES: O=S
Mol. weight [g/mol]: 48.06
CAS: 13827-32-2

Physical Properties

| Property code | Value | Unit | Source |
|---------------|--------------|--------|----------------|
| ea | 1.16 | eV | NIST Webbook |
| ea | 1.12 | eV | NIST Webbook |
| ea | 1.20 ± 0.10 | eV | NIST Webbook |
| ea | 1.10 | eV | NIST Webbook |
| ea | 1.09 ± 0.05 | eV | NIST Webbook |
| ea | 1.20 | eV | NIST Webbook |
| ea | 1.12 ± 0.01 | eV | NIST Webbook |
| ea | 1.10 | eV | NIST Webbook |
| gf | -163.83 | kJ/mol | Joback Method |
| hf | -137.45 | kJ/mol | Joback Method |
| hfus | 6.87 | kJ/mol | Joback Method |
| hvap | 28.03 | kJ/mol | Joback Method |
| ie | 10.34 ± 0.02 | eV | NIST Webbook |
| ie | 10.20 ± 0.03 | eV | NIST Webbook |
| ie | 10.00 ± 0.10 | eV | NIST Webbook |
| ie | 11.30 | eV | NIST Webbook |
| ie | 10.28 ± 0.02 | eV | NIST Webbook |
| ie | 10.31 | eV | NIST Webbook |
| ie | 10.29 ± 0.01 | eV | NIST Webbook |
| ie | 10.29 ± 0.00 | eV | NIST Webbook |
| ie | 10.29 ± 0.00 | eV | NIST Webbook |
| ie | 10.33 ± 0.01 | eV | NIST Webbook |
| log10ws | 0.66 | | Crippen Method |
| logp | -0.336 | | Crippen Method |
| mvol | 28.780 | ml/mol | McGowan Method |
| pc | 9174.85 | kPa | Joback Method |
| tb | 256.28 | K | Joback Method |
| tc | 421.10 | K | Joback Method |

| | | | |
|----|--------|----------------------|---------------|
| tf | 158.98 | K | Joback Method |
| vc | 0.107 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-------|---------|-----------------|---------------|
| cpg | 26.39 | J/mol×K | 256.28 | Joback Method |
| cpg | 28.49 | J/mol×K | 283.75 | Joback Method |
| cpg | 30.29 | J/mol×K | 311.22 | Joback Method |
| cpg | 31.81 | J/mol×K | 338.69 | Joback Method |
| cpg | 33.06 | J/mol×K | 366.16 | Joback Method |
| cpg | 34.07 | J/mol×K | 393.63 | Joback Method |
| cpg | 34.85 | J/mol×K | 421.10 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C13827322&Units=SI |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| ea: | Electron affinity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvap: | Enthalpy of vaporization at standard conditions |
| ie: | Ionization energy |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |

tb: Normal Boiling Point Temperature
tc: Critical Temperature
tf: Normal melting (fusion) point
vc: Critical Volume

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