

dinitrogen pentoxide

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| Other names: | dinitrogen pentoxide nitrogen oxide (N ₂ O ₅) nitrogen pentoxide |
| Inchi: | InChI=1S/N2O5/c3-1(4)7-2(5)6 |
| InchiKey: | ZWWCURLKEXEFQT-UHFFFAOYSA-N |
| Formula: | N ₂ O ₅ |
| SMILES: | O=[N+](O-)[N+](=O)[O-] |
| Mol. weight [g/mol]: | 108.01 |
| CAS: | 10102-03-1 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -84.78 | kJ/mol | Joback Method |
| hf | -197.07 | kJ/mol | Joback Method |
| hfus | 19.67 | kJ/mol | Joback Method |
| hvap | 51.19 | kJ/mol | Joback Method |
| ie | 12.30 | eV | NIST Webbook |
| ie | 11.40 | eV | NIST Webbook |
| log10ws | -1.06 | | Crippen Method |
| logp | -0.614 | | Crippen Method |
| mcvol | 51.570 | ml/mol | McGowan Method |
| pc | 6818.86 | kPa | Joback Method |
| tb | 525.50 | K | Joback Method |
| tc | 780.03 | K | Joback Method |
| tf | 399.21 | K | Joback Method |
| vc | 0.217 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 102.68 | J/mol×K | 525.50 | Joback Method |
| cpg | 106.56 | J/mol×K | 567.92 | Joback Method |
| cpg | 110.21 | J/mol×K | 610.34 | Joback Method |
| cpg | 113.60 | J/mol×K | 652.76 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 116.72 | J/mol×K | 695.18 | Joback Method |
| cpg | 119.57 | J/mol×K | 737.61 | Joback Method |
| cpg | 122.13 | J/mol×K | 780.03 | Joback Method |

Sources

| | |
|---|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C10102031&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Equilibrium Data for the N₂O₅ + HNO₃+ N₂O₄ System at 258.2 K, 265.2 K, 273.2 K and 298.2 K: | https://www.doi.org/10.1021/je800971f |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

Legend

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|-----------------|---|
| cpg: | Ideal gas heat capacity |
| 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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<https://www.chemeo.com/cid/13-201-3/dinitrogen-pentaoxide.pdf>

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