

Piperazine, 1,4-dinitro-

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| Other names: | N,N-Dinitropiperazine N,N'-Dinitropiperazine 1,4-Dinitropiperazine 1,4-Dinitro-1,4-diazacyclohexane |
| Inchi: | InChI=1S/C4H8N4O4/c9-7(10)5-1-2-6(4-3-5)8(11)12/h1-4H2 |
| InchiKey: | GSUMZAMXDRIYCL-UHFFFAOYSA-N |
| Formula: | C4H8N4O4 |
| SMILES: | O=[N+]([O-])N1CCN([N+](=O)[O-])CC1 |
| Mol. weight [g/mol]: | 176.13 |
| CAS: | 4164-37-8 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|-----------------|--------|----------------|
| chs | -2664.00 ± 2.00 | kJ/mol | NIST Webbook |
| hf | 58.00 ± 3.00 | kJ/mol | NIST Webbook |
| hfs | -53.00 ± 2.00 | kJ/mol | NIST Webbook |
| hsub | 111.00 ± 0.80 | kJ/mol | NIST Webbook |
| log10ws | -0.68 | | Crippen Method |
| logp | -1.013 | | Crippen Method |
| mvol | 111.160 | ml/mol | McGowan Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|---------------|--------|-----------------|--------------|
| hfust | 200.80 | kJ/mol | 489.20 | NIST Webbook |
| hfust | 33.93 | kJ/mol | 489.60 | NIST Webbook |
| hsubt | 111.00 ± 8.00 | kJ/mol | 342.50 | NIST Webbook |

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C4164378&Units=SI>

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307I>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>

Legend

chs: Standard solid enthalpy of combustion
hf: Enthalpy of formation at standard conditions
hfs: Solid phase enthalpy of formation at standard conditions
hfust: Enthalpy of fusion at a given temperature
hsub: Enthalpy of sublimation at standard conditions
hsubt: Enthalpy of sublimation at a given temperature
log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
mconvol: McGowan's characteristic volume

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