

Acetic acid, 4-nitrophenyl ester

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| Other names: | Acetic acid, p-nitrophenyl ester p-Nitrophenol acetate p-Nitrophenyl acetate 4-Nitrophenyl acetate p-Acetoxynitrobenzene p-Nitrobenzene acetate Phenol, p-nitro-, acetate NSC 2633 |
| Inchi: | InChI=1S/C8H7NO4/c1-6(10)13-8-4-2-7(3-5-8)9(11)12/h2-5H,1H3 |
| InchiKey: | QAUUDNIGJSLPSX-UHFFFAOYSA-N |
| Formula: | C8H7NO4 |
| SMILES: | CC(=O)Oc1ccc([N+](=O)[O-])cc1 |
| Mol. weight [g/mol]: | 181.15 |
| CAS: | 830-03-5 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|-------------|----------------------|----------------|
| gf | -79.11 | kJ/mol | Joback Method |
| hf | -238.95 | kJ/mol | Joback Method |
| hfus | 24.28 | kJ/mol | Joback Method |
| hvap | 62.09 | kJ/mol | Joback Method |
| ie | 9.50 ± 0.20 | eV | NIST Webbook |
| log10ws | -2.44 | | Crippen Method |
| logp | 1.520 | | Crippen Method |
| mcvol | 124.680 | ml/mol | McGowan Method |
| pc | 3848.31 | kPa | Joback Method |
| rinpola | 1430.00 | | NIST Webbook |
| tb | 642.23 | K | Joback Method |
| tc | 889.97 | K | Joback Method |
| tf | 434.63 | K | Joback Method |
| vc | 0.481 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 298.26 | J/mol×K | 642.23 | Joback Method |
| cpg | 308.88 | J/mol×K | 683.52 | Joback Method |
| cpg | 318.68 | J/mol×K | 724.81 | Joback Method |
| cpg | 327.67 | J/mol×K | 766.10 | Joback Method |
| cpg | 335.88 | J/mol×K | 807.39 | Joback Method |
| cpg | 343.31 | J/mol×K | 848.68 | Joback Method |
| cpg | 349.99 | J/mol×K | 889.97 | 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=C830035&Units=SI |

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 |
| rinpol: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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