

# Acetaldehyde, (2,4-dinitrophenyl)hydrazone

<b>Other names:</b>	Ethanal 2,4-dinitrophenylhydrazone NSC 403216 Ethanal DNPH, peak # 2
<b>Inchi:</b>	InChI=1S/C8H8N4O4/c1-2-9-10-7-4-3-6(11(13)14)5-8(7)12(15)16/h2-5,10H,1H3
<b>InchiKey:</b>	ONBOQRNOMHHDFB-UHFFFAOYSA-N
<b>Formula:</b>	C8H8N4O4
<b>SMILES:</b>	CC=NNc1ccc([N+](=O)[O-])cc1[N+](=O)[O-]
<b>Mol. weight [g/mol]:</b>	224.17
<b>CAS:</b>	1019-57-4

## Physical Properties

Property code	Value	Unit	Source
hf	119.31	kJ/mol	Joback Method
hvap	79.93	kJ/mol	Joback Method
log10ws	-3.36		Crippen Method
logp	1.921		Crippen Method
mcvol	150.320	ml/mol	McGowan Method
pc	3254.14	kPa	Joback Method
tb	849.61	K	Joback Method
tc	1123.10	K	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C1019574&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1019574&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>

## Legend

<b>hf:</b>	Enthalpy of formation at standard conditions
<b>h<sub>vap</sub>:</b>	Enthalpy of vaporization at standard conditions
<b>log<sub>10</sub>ws:</b>	Log <sub>10</sub> of Water solubility in mol/l
<b>log<sub>p</sub>:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

Latest version available from:

<https://www.cheméo.com/cid/14-512-7/Acetaldehyde-2-4-dinitrophenyl-hydrazone.pdf>

Generated by Cheméo on 2024-04-26 15:00:11.600002526 +0000 UTC m=+16432860.520579837.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.