

# Hydrazinecarboxamide, 2-(phenylmethylene)-

<b>Other names:</b>	Benzaldehyde, semicarbazone Benzylidenesemicarbazide 1-Benzylidenesemicarbazide
<b>Inchi:</b>	InChI=1S/C8H9N3O/c9-8(12)11-10-6-7-4-2-1-3-5-7/h1-6H,(H3,9,11,12)
<b>InchiKey:</b>	AKGUXECCGGCUDCV-UHFFFAOYSA-N
<b>Formula:</b>	C8H9N3O
<b>SMILES:</b>	NC(=O)NN=Cc1ccccc1
<b>Mol. weight [g/mol]:</b>	163.18
<b>CAS:</b>	1574-10-3

## Physical Properties

Property code	Value	Unit	Source
hf	84.98	kJ/mol	Joback Method
hvap	62.81	kJ/mol	Joback Method
log10ws	-1.91		Crippen Method
logp	0.689		Crippen Method
mcvol	127.030	ml/mol	McGowan Method
pc	3777.68	kPa	Joback Method
tb	662.37	K	Joback Method
tc	909.10	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=C1574103&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1574103&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>w<sub>s</sub>:</b>	Log <sub>10</sub> of Water solubility in mol/l
<b>log<sub>p</sub>:</b>	Octanol/Water partition coefficient
<b>mc<sub>vol</sub>:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

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