

# 2H-Indol-2-one, 1,3-dihydro-

<b>Other names:</b>	1,3-Dihydroindol-2-one 2,3-Dihydroindole-2-one 2,3-dihydroindol-2-one 2-Indolinone 2-Oxindole 2-Oxindoline 2-Oxoindoline Indol-2(3H)-one NSC 274863 Oxindol indolin-2-one o-(Aminophenyl)-acetic acid lactam oxindole
<b>Inchi:</b>	InChI=1S/C8H7NO/c10-8-5-6-3-1-2-4-7(6)9-8/h1-4H,5H2,(H,9,10)
<b>InchiKey:</b>	JYGFTBXVXVMTGB-UHFFFAOYSA-N
<b>Formula:</b>	C8H7NO
<b>SMILES:</b>	O=C1Cc2ccccc2N1
<b>Mol. weight [g/mol]:</b>	133.15
<b>CAS:</b>	59-48-3

## Physical Properties

Property code	Value	Unit	Source
gf	152.84	kJ/mol	Joback Method
hf	9.86	kJ/mol	Joback Method
hfus	16.29	kJ/mol	Joback Method
hvap	47.57	kJ/mol	Joback Method
ie	8.36	eV	NIST Webbook
log10ws	-1.53		Crippen Method
logp	1.181		Crippen Method
mcvol	100.510	ml/mol	McGowan Method
pc	4775.98	kPa	Joback Method
tb	541.88	K	Joback Method
tc	796.73	K	Joback Method
tf	414.29	K	Joback Method
tt	393.93	K	Solubility Measurement and Correlation of 2-Oxindole in 12 Pure Organic Solvents

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	218.51	J/mol×K	541.88	Joback Method
cpg	230.99	J/mol×K	584.36	Joback Method
cpg	242.64	J/mol×K	626.83	Joback Method
cpg	253.49	J/mol×K	669.31	Joback Method
cpg	263.56	J/mol×K	711.78	Joback Method
cpg	272.90	J/mol×K	754.26	Joback Method
cpg	281.52	J/mol×K	796.73	Joback Method

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	500.20	K	9.70	NIST Webbook
tbrp	500.20	K	3.10	NIST Webbook

## Sources

Solubility Measurement and Correlation of 2-Oxindole in 12 Pure Organic Solvents	<a href="https://www.doi.org/10.1021/acs.jced.9b00308">https://www.doi.org/10.1021/acs.jced.9b00308</a>
Joback Method:	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
McGowan Method:	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
NIST Webbook:	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C59483&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C59483&amp;Units=SI</a>
Crippen Method:	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
Crippen Method:	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>

## Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation

<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>h vap:</b>	Enthalpy of vaporization at standard conditions
<b>ie:</b>	Ionization energy
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</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>tbrp:</b>	Boiling point at reduced pressure
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>tt:</b>	Triple Point Temperature
<b>vc:</b>	Critical Volume

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