

# Benzofurazan, 4-nitro-, 1-oxide

<b>Other names:</b>	4-Nitrobenzofuroxan B2789 Nitrobenzofuroxan 2,1,3-Benzoxadiazole, 4-nitro-, 1-oxide 4-Nitrobenzofuroxan-1-oxide
<b>Inchi:</b>	InChI=1S/C6H3N3O4/c10-8(11)4-2-1-3-5-6(4)7-13-9(5)12/h1-3H
<b>InchiKey:</b>	XUCKCDGEIRMZPM-UHFFFAOYSA-N
<b>Formula:</b>	C6H3N3O4
<b>SMILES:</b>	O=[N+](O-)c1cccc2c1no[n+]2[O-]
<b>Mol. weight [g/mol]:</b>	181.11
<b>CAS:</b>	18771-85-2

## Physical Properties

Property code	Value	Unit	Source
chs	-2992.40 ± 1.90	kJ/mol	NIST Webbook
hf	299.90 ± 2.60	kJ/mol	NIST Webbook
hfs	202.60 ± 2.10	kJ/mol	NIST Webbook
hsub	97.30 ± 1.60	kJ/mol	NIST Webbook
hsub	97.30 ± 1.60	kJ/mol	NIST Webbook
log10ws	-8.95		Crippen Method
logp	0.369		Crippen Method
mvol	105.600	ml/mol	McGowan Method

## Sources

<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=C18771852&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C18771852&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>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>

# Legend

<b>chs:</b>	Standard solid enthalpy of combustion
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfs:</b>	Solid phase enthalpy of formation at standard conditions
<b>hsub:</b>	Enthalpy of sublimation at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume

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