

# 9H-Fluoren-9-one, 2,7-dinitro-

<b>Other names:</b>	2,7-Dinitrofluoren-9-one 2,7-Dinitro-9-fluorenone Fluoren-9-one, 2,7-dinitro- 2,7-Dinitrofluorenone 9-Fluorenone, 2,7-dinitro- 2,7-Dinitro-9H-fluoren-9-one
<b>Inchi:</b>	InChI=1S/C13H6N2O5/c16-13-11-5-7(14(17)18)1-3-9(11)10-4-2-8(15(19)20)6-12(10)13/
<b>InchiKey:</b>	HDVGAFBXTXDYIB-UHFFFAOYSA-N
<b>Formula:</b>	C13H6N2O5
<b>SMILES:</b>	O=C1c2cc([N+](=O)[O-])ccc2-c2ccc([N+](=O)[O-])cc21
<b>Mol. weight [g/mol]:</b>	270.20
<b>CAS:</b>	31551-45-8

## Physical Properties

Property code	Value	Unit	Source
gf	286.05	kJ/mol	Joback Method
hf	61.77	kJ/mol	Joback Method
hfus	39.45	kJ/mol	Joback Method
hvap	89.04	kJ/mol	Joback Method
log10ws	-5.58		Crippen Method
logp	2.714		Crippen Method
mcvol	172.060	ml/mol	McGowan Method
pc	3493.01	kPa	Joback Method
rinpol	432.66		NIST Webbook
tb	944.49	K	Joback Method
tc	1237.67	K	Joback Method
tf	723.85	K	Joback Method
vc	0.693	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	491.54	J/molxK	944.49	Joback Method
cpg	500.72	J/molxK	993.35	Joback Method

cpg	509.24	J/mol×K	1042.22	Joback Method
cpg	517.25	J/mol×K	1091.08	Joback Method
cpg	524.90	J/mol×K	1139.94	Joback Method
cpg	532.34	J/mol×K	1188.80	Joback Method
cpg	539.71	J/mol×K	1237.67	Joback 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=C31551458&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C31551458&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>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>gf:</b>	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>hvap:</b>	Enthalpy of vaporization 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
<b>pc:</b>	Critical Pressure
<b>rinpola:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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