

# 9H-Fluoren-9-ol

<b>Other names:</b>	9-fluorenol 9-hydroxyfluorene 9H-Fluorene-9-ol Diphenylene carbinol Fluoren-9-ol Fluorenol
<b>Inchi:</b>	InChI=1S/C13H10O/c14-13-11-7-3-1-5-9(11)10-6-2-4-8-12(10)13/h1-8,13-14H
<b>InchiKey:</b>	AFMVESZYOYKHDBJ-UHFFFAOYSA-N
<b>Formula:</b>	C13H10O
<b>SMILES:</b>	OC1c2ccccc2-c2ccccc21
<b>Mol. weight [g/mol]:</b>	182.22
<b>CAS:</b>	1689-64-1

## Physical Properties

Property code	Value	Unit	Source
gf	212.27	kJ/mol	Joback Method
hf	71.36	kJ/mol	Joback Method
hfus	24.20	kJ/mol	Experimental and Computational Study of the Thermodynamic Properties of 9-Fluorenone and 9-Fluorenol
hvap	66.66	kJ/mol	Joback Method
log10ws	-4.14		Crippen Method
logp	2.749		Crippen Method
mcvol	141.520	ml/mol	McGowan Method
pc	3620.24	kPa	Joback Method
rinpol	1791.00		NIST Webbook
rinpol	1704.00		NIST Webbook
rinpol	280.50		NIST Webbook
rinpol	280.50		NIST Webbook
rinpol	1791.00		NIST Webbook
ripol	2473.00		NIST Webbook
tb	650.54	K	Joback Method
tc	876.67	K	Joback Method
tf	427.60 ± 2.00	K	NIST Webbook
vc	0.539	m <sup>3</sup> /kmol	Joback Method

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	357.87	J/molxK	650.54	Joback Method
cpg	369.52	J/molxK	688.23	Joback Method
cpg	380.30	J/molxK	725.92	Joback Method
cpg	390.30	J/molxK	763.60	Joback Method
cpg	399.63	J/molxK	801.29	Joback Method
cpg	408.38	J/molxK	838.98	Joback Method
cpg	416.67	J/molxK	876.67	Joback Method
dvisc	0.0013355	Paxs	441.72	Joback Method
dvisc	0.0022872	Paxs	399.95	Joback Method
dvisc	0.0008557	Paxs	483.48	Joback Method
dvisc	0.0005886	Paxs	525.25	Joback Method
dvisc	0.0004277	Paxs	567.01	Joback Method
dvisc	0.0003248	Paxs	608.77	Joback Method
dvisc	0.0002555	Paxs	650.54	Joback Method
hvapt	50.40	kJ/mol	435.00	NIST Webbook
hvapt	49.70	kJ/mol	465.00	NIST Webbook
hvapt	48.90	kJ/mol	505.00	NIST Webbook
psub	2.10e-05	kPa	326.10	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	2.44e-05	kPa	327.50	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique

psub	3.15e-05	kPa	329.80	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	3.74e-05	kPa	331.30	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	5.12e-05	kPa	334.90	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	1.01e-04	kPa	340.90	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	1.15e-04	kPa	342.60	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique

psub	1.23e-04	kPa	343.60	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	1.44e-04	kPa	344.70	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	1.59e-04	kPa	346.10	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	1.86e-04	kPa	347.40	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	2.76e-04	kPa	351.20	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique

psub	4.02e-04	kPa	355.10	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	5.60e-04	kPa	358.70	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	7.40e-04	kPa	361.90	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique
psub	1.03e-03	kPa	366.10	Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique

## Sources

### Crippen Method:

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

Vapor pressures and sublimation enthalpies of seven heteroatomic aromatic hydrocarbons measured using the Knudsen effusion technique: Joback Method and 9-Fluoreno1 :

<https://www.doi.org/10.1016/j.jct.2010.01.014>

Experimental and Computational Study of the Thermodynamic Properties of 9-Fluoreno1 and 9-Fluoreno1 :

<https://www.doi.org/10.1021/je300584m>

### McGowan Method:

[https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

<http://link.springer.com/article/10.1007/BF02311772>

### NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C1689641&Units=SI>

### Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

# Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<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>hvapt:</b>	Enthalpy of vaporization at a given temperature
<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>psub:</b>	Sublimation pressure
<b>rinpol:</b>	Non-polar retention indices
<b>ripol:</b>	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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