

# 1-Nonene, 1,1,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,9-octadecafluoro-

Other names:

perfluoronon-1-ene

perfluoro-1-nonene

Inchi: InChI=1S/C9F18/c10-1(2(11)12)3(13,14)4(15,16)5(17,18)6(19,20)7(21,22)8(23,24)9(25,26)

InchiKey: UAFOIVDGAVVKTE-UHFFFAOYSA-N

Formula: C9F18

SMILES: FC(F)=C(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)F

Mol. weight [g/mol]: 450.07

CAS: 376-22-7

## Physical Properties

Property code	Value	Unit	Source
gf	-3398.68	kJ/mol	Joback Method
hf	-3722.68	kJ/mol	Joback Method
hfus	20.19	kJ/mol	Joback Method
hvap	11.97	kJ/mol	Joback Method
log10ws	-7.04		Crippen Method
logp	6.438		Crippen Method
mcvol	165.230	ml/mol	McGowan Method
pc	1374.80	kPa	Joback Method
rinpola	423.00		NIST Webbook
tb	396.00 ± 3.00	K	NIST Webbook
tc	483.99	K	Joback Method
tf	185.75	K	Joback Method
vc	0.768	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	408.73	J/mol×K	373.49	Joback Method
cpg	422.19	J/mol×K	391.91	Joback Method
cpg	434.86	J/mol×K	410.32	Joback Method
cpg	446.77	J/mol×K	428.74	Joback Method
cpg	457.93	J/mol×K	447.16	Joback Method
cpg	468.39	J/mol×K	465.57	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=C376227&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C376227&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</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>rinpol:</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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