

# Benzene, nonyl-

<b>Other names:</b>	1-PHENYLNONANE NONYLBENZENE Nonane, 1-phenyl- n-Nonylbenzene
<b>Inchi:</b>	InChI=1S/C15H24/c1-2-3-4-5-6-7-9-12-15-13-10-8-11-14-15/h8,10-11,13-14H,2-7,9,12H2
<b>InchiKey:</b>	LIXVMPBOGDCSRM-UHFFFAOYSA-N
<b>Formula:</b>	C15H24
<b>SMILES:</b>	CCCCCCCCCc1ccccc1
<b>Mol. weight [g/mol]:</b>	204.35
<b>CAS:</b>	1081-77-2

## Physical Properties

Property code	Value	Unit	Source
af	0.6280		KDB
gf	187.83	kJ/mol	Joback Method
hf	-116.40	kJ/mol	Joback Method
hfus	28.65	kJ/mol	Joback Method
hvap	74.80	kJ/mol	NIST Webbook
hvap	74.10 ± 0.50	kJ/mol	NIST Webbook
log10ws	-5.20		Crippen Method
logp	4.980		Crippen Method
mcpol	198.450	ml/mol	McGowan Method
nfpaf	%!d(float64=1)		KDB
pc	1900.00	kPa	KDB
rinpol	1527.00		NIST Webbook
rinpol	1555.00		NIST Webbook
rinpol	1586.00		NIST Webbook
rinpol	1559.50		NIST Webbook
rinpol	1560.00		NIST Webbook
rinpol	1552.20		NIST Webbook
rinpol	1556.98		NIST Webbook
rinpol	1560.39		NIST Webbook
rinpol	1570.95		NIST Webbook
rinpol	1576.27		NIST Webbook
rinpol	1579.75		NIST Webbook
rinpol	1527.00		NIST Webbook
rinpol	1554.10		NIST Webbook

rinpol	1542.80		NIST Webbook
rinpol	1554.00		NIST Webbook
rinpol	1542.80		NIST Webbook
rinpol	1555.00		NIST Webbook
rinpol	1527.00		NIST Webbook
rinpol	265.60		NIST Webbook
rinpol	265.51		NIST Webbook
rinpol	265.60		NIST Webbook
rinpol	1555.00		NIST Webbook
rinpol	1555.00		NIST Webbook
rinpol	1554.10		NIST Webbook
ripol	1854.00		NIST Webbook
ripol	1821.00		NIST Webbook
ripol	1854.00		NIST Webbook
ripol	1847.00		NIST Webbook
tb	555.20	K	KDB
tc	741.00	K	KDB
tf	249.00	K	KDB
vc	0.790	m <sup>3</sup> /kmol	KDB
zc	0.2436270		KDB

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	576.51	J/molxK	728.92	Joback Method
cpg	561.51	J/molxK	696.99	Joback Method
cpg	545.65	J/molxK	665.07	Joback Method
cpg	528.89	J/molxK	633.14	Joback Method
cpg	511.21	J/molxK	601.21	Joback Method
cpg	492.55	J/molxK	569.28	Joback Method
cpg	590.69	J/molxK	760.85	Joback Method
dvisc	0.0035750	Paxs	285.23	Joback Method
dvisc	0.0001656	Paxs	569.28	Joback Method
dvisc	0.0002190	Paxs	521.94	Joback Method
dvisc	0.0003064	Paxs	474.60	Joback Method
dvisc	0.0004617	Paxs	427.25	Joback Method
dvisc	0.0007706	Paxs	379.91	Joback Method
dvisc	0.0014881	Paxs	332.57	Joback Method
hvapt	48.95	kJ/mol	555.20	KDB
hvapt	69.70	kJ/mol	365.50	NIST Webbook

# Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.54330e+01
Coeff. B	-5.20756e+03
Coeff. C	-7.04730e+01
Temperature range (K), min.	414.31
Temperature range (K), max.	584.98

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/T + C \cdot \ln(T) + D \cdot T^2$
Coeff. A	1.14515e+02
Coeff. B	-1.27354e+04
Coeff. C	-1.39957e+01
Coeff. D	4.83248e-06
Temperature range (K), min.	249.00
Temperature range (K), max.	741.00

# Sources

<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>KDB:</b>	<a href="https://www.thermo.com/files/research/kdb/mol/mol711.mol">https://www.thermo.com/files/research/kdb/mol/mol711.mol</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=C1081772&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1081772&amp;Units=SI</a>
<b>The Yaws Handbook of Vapor Pressure:</b>	<a href="https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure">https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure</a>
<b>KDB Vapor Pressure Data:</b>	<a href="https://www.thermo.com/research/kdb/hcprop/showprop.php?cmpid=711">https://www.thermo.com/research/kdb/hcprop/showprop.php?cmpid=711</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.chemo.com/doc/models/crippen_log10ws">https://www.chemo.com/doc/models/crippen_log10ws</a>

# Legend

<b>af:</b>	Acentric Factor
<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>nfpa:</b>	NFPA Fire Rating
<b>pc:</b>	Critical Pressure
<b>pvap:</b>	Vapor 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
<b>zc:</b>	Critical Compressibility

Latest version available from:

<https://www.cheméo.com/cid/20-664-2/Benzene-nonyl.pdf>

Generated by Cheméo on 2024-04-24 09:51:41.061036929 +0000 UTC m=+16241549.981614245.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.