

# Hydrazine, 1-butyl-1-methyl-

<b>Other names:</b>	1-Butyl-1-methylhydrazine 1-Methyl-1-n-butylhydrazine Hydrazine, butylmethyl- Methylbutyl hydrazine
<b>Inchi:</b>	InChI=1S/C5H14N2/c1-3-4-5-7(2)6/h3-6H2,1-2H3
<b>InchiKey:</b>	OEEGQMYUIKPQMM-UHFFFAOYSA-N
<b>Formula:</b>	C5H14N2
<b>SMILES:</b>	CCCCN(C)N
<b>Mol. weight [g/mol]:</b>	102.18
<b>CAS:</b>	20240-62-4

## Physical Properties

Property code	Value	Unit	Source
gf	168.45	kJ/mol	Joback Method
hf	-45.21	kJ/mol	Joback Method
hfus	16.92	kJ/mol	Joback Method
hvap	39.41	kJ/mol	Joback Method
ie	7.62 ± 0.05	eV	NIST Webbook
ie	7.51 ± 0.02	eV	NIST Webbook
ie	7.82 ± 0.05	eV	NIST Webbook
log10ws	-0.91		Crippen Method
logp	0.592		Crippen Method
mcvol	101.270	ml/mol	McGowan Method
pc	3704.46	kPa	Joback Method
rinpola	798.00		NIST Webbook
tb	398.77	K	Joback Method
tc	578.11	K	Joback Method
tf	261.84	K	Joback Method
vc	0.362	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	201.04	J/mol×K	398.77	Joback Method

cpg	212.39	J/mol×K	428.66	Joback Method
cpg	223.25	J/mol×K	458.55	Joback Method
cpg	233.62	J/mol×K	488.44	Joback Method
cpg	243.52	J/mol×K	518.33	Joback Method
cpg	252.96	J/mol×K	548.22	Joback Method
cpg	261.97	J/mol×K	578.11	Joback Method

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	2.07602e+01
Coeff. B	-5.35625e+03
Coeff. C	-5.87300e+01
Temperature range (K), min.	320.36
Temperature range (K), max.	405.44

## Sources

<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>
<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=C20240624&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C20240624&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>

## 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>ie:</b>	Ionization energy

<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>pvap:</b>	Vapor 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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