

# nitric acid

Other names:	AZOTIC ACID HYDROGEN NITRATE
Inchi:	InChI=1S/HNO3/c2-1(3)4/h(H,2,3,4)
InchiKey:	GRYLNZFGIOXLOG-UHFFFAOYSA-N
Formula:	HNO <sub>3</sub>
SMILES:	O=[N+](O-)O
Mol. weight [g/mol]:	63.01
CAS:	7697-37-2

## Physical Properties

Property code	Value	Unit	Source
affp	751.40	kJ/mol	NIST Webbook
affp	670.30	kJ/mol	NIST Webbook
basg	731.50	kJ/mol	NIST Webbook
ea	0.57 ± 0.15	eV	NIST Webbook
ea	0.56 ± 0.17	eV	NIST Webbook
gf	-152.15	kJ/mol	Joback Method
hf	-206.32	kJ/mol	Joback Method
hfus	11.21	kJ/mol	Joback Method
hvap	48.86	kJ/mol	Joback Method
ie	11.03 ± 0.01	eV	NIST Webbook
ie	12.20	eV	NIST Webbook
ie	11.96	eV	NIST Webbook
ie	11.95 ± 0.01	eV	NIST Webbook
log10ws	0.44		Crippen Method
logp	-0.348		Crippen Method
mcvol	34.150	ml/mol	McGowan Method
nfpah	%!d(float64=2)		KDB
nfpas	%!d(float64=1)		KDB
pc	8058.04	kPa	Joback Method
tb	443.42	K	Joback Method
tc	648.46	K	Joback Method
tf	294.19	K	Joback Method
vc	0.137	m <sup>3</sup> /kmol	Joback Method

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	59.38	J/mol×K	443.42	Joback Method
cpg	61.79	J/mol×K	477.59	Joback Method
cpg	64.07	J/mol×K	511.77	Joback Method
cpg	66.21	J/mol×K	545.94	Joback Method
cpg	68.23	J/mol×K	580.12	Joback Method
cpg	70.12	J/mol×K	614.29	Joback Method
cpg	71.89	J/mol×K	648.46	Joback Method
hvapt	38.60	kJ/mol	314.50	NIST Webbook

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.54030e+01
Coeff. B	-3.29672e+03
Coeff. C	-5.04600e+01
Temperature range (K), min.	231.55
Temperature range (K), max.	376.10

## Sources

### NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C7697372&Units=SI>

### Thermodynamic Properties of Ternary Solutions in the Water-Nitric Acid-Rare Earth Nitrate Equilibrium Systems at Nitrobenzene-Inorganic Acid Systems at 298.15 K: Joback Method:

<https://www.doi.org/10.1021/acs.jced.6b00357>

<https://www.doi.org/10.1021/je800878b>

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

### The Yaws Handbook of Vapor Pressure: Crippen Method:

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

### Thermochemical Analysis on Rare Earth Complex of Gadolinium with Salicylic acid and 8-hydroxyquinoline:

<https://www.doi.org/10.1016/j.tca.2012.08.027>

<https://www.chemistry.org/research/kdb/hcprop/showprop.php?cmpid=1911>

### Apparent molar volumes and apparent molar heat capacities of aqueous lead(II) nitrate, Mn(II) nitrate, and Cu(II) nitrate in the N2O5-HNO3-H2O system at 268.2 K, 265.0 K, 273.2 K, and 281.2 K in liquid equilibrium: Nitric acid - water - anisole/4-methyl anisole:

<https://www.doi.org/10.1016/j.jct.2003.09.013>

<https://www.doi.org/10.1021/je800971f>

<https://www.doi.org/10.1016/j.fluid.2011.10.022>

Densities and Excess Molar Volumes of the Ternary System N <sub>2</sub> O <sub>4</sub> + H <sub>2</sub> O + HF at 273 K, 283 K, 289 K, tetrahydrofuran on the phase equilibrium of systems containing 5-hydroxymethylfurfural, water, organic solvent in the absence and presence of sodium chloride:	<a href="https://www.doi.org/10.1021/je101357s">https://www.doi.org/10.1021/je101357s</a>
Cripen Method:	<a href="https://www.doi.org/10.1016/j.jct.2019.01.001">https://www.doi.org/10.1016/j.jct.2019.01.001</a>
Equilibrium Methods:	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
McCowan Method:	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
Distillation Separation of Hydrofluoric Acid and Nitric Acid From Acid Waste using the Salt Effect on Vapor-Liquid Equilibrium :	<a href="https://www.doi.org/10.1007/s10765-010-0904-8">https://www.doi.org/10.1007/s10765-010-0904-8</a>

## Legend

<b>affp:</b>	Proton affinity
<b>basg:</b>	Gas basicity
<b>cpg:</b>	Ideal gas heat capacity
<b>ea:</b>	Electron affinity
<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>ie:</b>	Ionization energy
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McCowan's characteristic volume
<b>nfpah:</b>	NFPA Health Rating
<b>nfpas:</b>	NFPA Safety Rating
<b>pc:</b>	Critical Pressure
<b>pvap:</b>	Vapor pressure
<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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