

# Nitrous oxide

<b>Other names:</b>	DINITROGEN MONOXIDE Dinitrogen oxide Factitious air Hyponitrous acid anhydride LAUGHING GAS N2O Nitral Nitrogen monoxide Nitrogen oxide Nitrogen oxide (N2O) UN 1070 UN 2201
<b>Inchi:</b>	InChI=1S/N2O/c1-2-3
<b>InchiKey:</b>	GQPLMRYTRLFLPF-UHFFFAOYSA-N
<b>Formula:</b>	N2O
<b>SMILES:</b>	[N-]=[N+]=O
<b>Mol. weight [g/mol]:</b>	44.01
<b>CAS:</b>	10024-97-2

## Physical Properties

Property code	Value	Unit	Source
af	0.1650		KDB
affp	575.20	kJ/mol	NIST Webbook
affp	549.80	kJ/mol	NIST Webbook
basg	548.70	kJ/mol	NIST Webbook
basg	523.30	kJ/mol	NIST Webbook
dm	0.20	debye	KDB
ea	0.27 ± 0.17	eV	NIST Webbook
ea	0.76 ± 0.10	eV	NIST Webbook
ea	0.22 ± 0.10	eV	NIST Webbook
ea	-0.15 ± 0.10	eV	NIST Webbook
gf	103.70	kJ/mol	KDB
hf	81.60	kJ/mol	KDB
ie	12.89	eV	NIST Webbook
ie	12.80 ± 0.05	eV	NIST Webbook
ie	12.88 ± 0.01	eV	NIST Webbook
ie	12.89 ± 0.00	eV	NIST Webbook

ie	12.89	eV	NIST Webbook
ie	12.00 ± 1.00	eV	NIST Webbook
ie	12.91 ± 0.03	eV	NIST Webbook
ie	12.89 ± 0.00	eV	NIST Webbook
ie	12.88 ± 0.01	eV	NIST Webbook
ie	12.89 ± 0.01	eV	NIST Webbook
ie	12.90	eV	NIST Webbook
ie	12.89 ± 0.01	eV	NIST Webbook
ie	12.89 ± 0.01	eV	NIST Webbook
ie	12.89	eV	NIST Webbook
ie	12.89 ± 0.01	eV	NIST Webbook
ie	12.89	eV	NIST Webbook
log10ws	1.67		Crippen Method
logp	-0.158		Crippen Method
mcvol	28.090	ml/mol	McGowan Method
pc	7255.00	kPa	KDB
pc	7270.00 ± 50.00	kPa	NIST Webbook
pc	7234.60 ± 50.66	kPa	NIST Webbook
pc	7238.00 ± 20.00	kPa	NIST Webbook
pt	87.89 ± 0.01	kPa	NIST Webbook
pt	87.91 ± 0.01	kPa	NIST Webbook
pt	87.85	kPa	KDB
rhoc	453.33 ± 4.40	kg/m3	NIST Webbook
rhoc	448.93	kg/m3	NIST Webbook
rhoc	453.33 ± 2.20	kg/m3	NIST Webbook
rinpol	182.00		NIST Webbook
rinpol	182.00		NIST Webbook
tb	184.67	K	KDB
tc	309.57	K	KDB
tc	309.56 ± 0.15	K	NIST Webbook
tc	309.65 ± 0.20	K	NIST Webbook
tc	309.49	K	NIST Webbook
tc	309.55 ± 0.50	K	NIST Webbook
tf	182.30	K	KDB
tt	182.32	K	KDB
tt	181.99	K	Solid Liquid Equilibria for the CO2 + R23 and N2O + R23 Systems
vc	0.097	m3/kmol	KDB
vc	0.096 ± 0.002	m3/kmol	NIST Webbook
zc	0.2734100		KDB
zra	0.28		KDB

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hfust	6.50	kJ/mol	182.40	NIST Webbook
hsubt	23.60	kJ/mol	113.00	NIST Webbook
hsubt	24.60	kJ/mol	165.00	NIST Webbook
hsubt	25.10 ± 0.40	kJ/mol	74.00	NIST Webbook
hvapt	16.10	kJ/mol	209.00	NIST Webbook
hvapt	16.50	kJ/mol	184.70	NIST Webbook
pvap	2065.40	kPa	257.87	PVT Properties of Dinitrogen Monoxide
pvap	1092.90	kPa	237.49	PVT Properties of Dinitrogen Monoxide
pvap	1294.70	kPa	242.57	PVT Properties of Dinitrogen Monoxide
pvap	1521.90	kPa	247.65	PVT Properties of Dinitrogen Monoxide
pvap	1777.00	kPa	252.73	PVT Properties of Dinitrogen Monoxide
pvap	2061.80	kPa	257.80	PVT Properties of Dinitrogen Monoxide
pvap	915.40	kPa	232.43	PVT Properties of Dinitrogen Monoxide
pvap	2379.00	kPa	262.90	PVT Properties of Dinitrogen Monoxide
pvap	2381.80	kPa	262.95	PVT Properties of Dinitrogen Monoxide
pvap	2730.60	kPa	267.98	PVT Properties of Dinitrogen Monoxide
pvap	2734.00	kPa	268.03	PVT Properties of Dinitrogen Monoxide
pvap	3119.90	kPa	273.08	PVT Properties of Dinitrogen Monoxide
pvap	912.10	kPa	232.37	PVT Properties of Dinitrogen Monoxide
pvap	760.60	kPa	227.36	PVT Properties of Dinitrogen Monoxide

pvap	759.60	kPa	227.32	PVT Properties of Dinitrogen Monoxide
pvap	652.80	kPa	223.31	PVT Properties of Dinitrogen Monoxide
pvap	626.30	kPa	222.28	PVT Properties of Dinitrogen Monoxide
pvap	554.70	kPa	219.24	PVT Properties of Dinitrogen Monoxide
rho1	1226.00	kg/m3	184.00	KDB

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.36769e+01
Coeff. B	-1.26781e+03
Coeff. C	-4.47100e+01
Temperature range (K), min.	182.30
Temperature range (K), max.	309.57

## Sources

Density, Viscosity and N<sub>2</sub>O solubility of Aqueous 2- (methylamino)ethanol  
 Solubility of N<sub>2</sub>O in and Density and Viscosity of Aqueous Solutions of Physical Solubility and Diffusivity of N<sub>2</sub>O and CO<sub>2</sub> in Aqueous Sodium Bicarbonate Solutions, and N<sub>2</sub>O Solubility of Aqueous Solutions of MEA, Solid-Liquid Equilibria for the CO<sub>2</sub>+R<sub>15</sub> and R<sub>15</sub>+R<sub>15</sub> Systems: An experimental and modeling study of physical N<sub>2</sub>O solubility in 2-(methylamino)ethanol + (1-x)HCl at temperatures (159.01 and 182.3) K  
 Binary Diffusion Coefficient Data of Various Gas Systems Determined  
 Solubility of carbon dioxide, nitrous oxide, dinitrogen and nitrogen in 1-butyl-3-methylpyridinium and hexadecyltetradecylphosphonium mesitylfluorobenzenesulfonate ionic liquids  
 Physical and chemical properties of Novel ionic liquids based on propylamine and diethylamine  
 Isochores PVT measurements for the C<sub>2</sub>H<sub>6</sub> + N<sub>2</sub>O binary system: Phase Behavior of N<sub>2</sub>O and CO<sub>2</sub> in Room-Temperature Ionic Liquids [bmim][Tf<sub>2</sub>N], [bmim][BF<sub>4</sub>], [bmim][N(CN)<sub>2</sub>], [bmim][Ac], [eam][NO<sub>3</sub>], and [bmim][SCN]:

<https://www.doi.org/10.1021/acs.jced.6b00504>  
<https://www.doi.org/10.1021/je300102d>  
<https://www.doi.org/10.1021/je0503913>  
<https://www.doi.org/10.1021/acs.jced.8b00070>  
<https://www.doi.org/10.1021/je700384e>  
<https://www.doi.org/10.1016/j.jct.2019.06.008>  
<https://www.doi.org/10.1016/j.jct.2004.05.009>  
<https://www.doi.org/10.1007/s10765-015-1981-5>  
<https://www.doi.org/10.1016/j.jct.2012.11.010>  
<https://www.doi.org/10.1021/je049842u>  
<https://www.doi.org/10.1021/acs.jced.5b00922>  
<https://www.doi.org/10.1021/je800409d>  
<https://www.doi.org/10.1021/je0503148>  
<https://www.doi.org/10.1007/s10765-011-1150-4>

<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>Solid-Liquid Equilibria for the CO<sub>2</sub> + R125 and N<sub>2</sub>O + R125 Systems: A New Solid-Liquid Equilibria for the CO<sub>2</sub> + R23 and N<sub>2</sub>O + R23 Systems:</b>	<a href="https://www.doi.org/10.1021/je0603067">https://www.doi.org/10.1021/je0603067</a> <a href="https://www.doi.org/10.1007/s10765-008-0511-0">https://www.doi.org/10.1007/s10765-008-0511-0</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>Vapour pressure and excess Gibbs free energy of the ternary system (x1CH<sub>3</sub>F, x2N<sub>2</sub>O, x3CO<sub>2</sub>) at a temperature of 298.15 K:</b>	<a href="https://www.doi.org/10.1016/j.jct.2006.04.010">https://www.doi.org/10.1016/j.jct.2006.04.010</a> <a href="https://www.doi.org/10.1021/je049633+">https://www.doi.org/10.1021/je049633+</a>
<b>Solubility of Nitrous Oxide in Aqueous Methyldiethanolamine Solutions:</b>	<a href="https://www.doi.org/10.1021/acs.jced.7b00112">https://www.doi.org/10.1021/acs.jced.7b00112</a>
<b>Experimental Study of a Hydrophobic Solvent for Natural Gas Sweetening Based on the Solubility and Selectivity of Light Hydrocarbons (CH<sub>4</sub>, C<sub>2</sub>H<sub>6</sub>) and Acid Gases (CO<sub>2</sub> and H<sub>2</sub>S) at 298-353 K:</b>	<a href="https://www.doi.org/10.1021/acs.jced.8b00735">https://www.doi.org/10.1021/acs.jced.8b00735</a> <a href="https://www.doi.org/10.1021/je800460a">https://www.doi.org/10.1021/je800460a</a> <a href="https://www.cheric.org/files/research/kdb/mol/mol1931.mol">https://www.cheric.org/files/research/kdb/mol/mol1931.mol</a>
<b>Physicochemical Properties of Aqueous Solutions of 2-(Diethylamino)-ethanol:</b>	<a href="https://www.doi.org/10.1021/je9007926">https://www.doi.org/10.1021/je9007926</a>
<b>Burnett Measurements and Virial Coefficients for the R32+N<sub>2</sub>O System: PVTx measurements for N<sub>2</sub>O+CH<sub>3</sub>F, N<sub>2</sub>O+CH<sub>2</sub>F<sub>2</sub>, and N<sub>2</sub>O+ CHF<sub>3</sub> binary systems:</b>	<a href="https://www.doi.org/10.1007/s10765-006-0033-6">https://www.doi.org/10.1007/s10765-006-0033-6</a> <a href="https://www.doi.org/10.1016/j.fluid.2004.11.021">https://www.doi.org/10.1016/j.fluid.2004.11.021</a> <a href="https://www.doi.org/10.1007/s10765-006-0082-x">https://www.doi.org/10.1007/s10765-006-0082-x</a>
<b>Virial Coefficients from Burnett Measurements for the R23 + N<sub>2</sub>O system:</b>	<a href="https://www.doi.org/10.1016/j.fluid.2008.04.003">https://www.doi.org/10.1016/j.fluid.2008.04.003</a>
<b>Solubility of carbon dioxide in aqueous solutions of isomeric PVTx measurements for the N<sub>2</sub>O + R125 Binary System:</b>	<a href="https://www.doi.org/10.1021/je060140+">https://www.doi.org/10.1021/je060140+</a>
<b>Solubility of N<sub>2</sub>O and CO<sub>2</sub> in non-aqueous systems of 1,4-Butanediamine and Vapor Pressure Measurements and model correlations for carbon dioxide, nitrous oxide and methane in ionic liquids at pressures close to atmospheric:</b>	<a href="https://www.doi.org/10.1016/j.jct.2019.05.017">https://www.doi.org/10.1016/j.jct.2019.05.017</a> <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> <a href="https://www.doi.org/10.1016/j.fluid.2014.03.015">https://www.doi.org/10.1016/j.fluid.2014.03.015</a> <a href="https://www.doi.org/10.1021/acs.jced.6b00013">https://www.doi.org/10.1021/acs.jced.6b00013</a>
<b>Physicochemical Properties of Aqueous Potassium Salts of Basic Amino Acids and Correlations of Solubility of N<sub>2</sub>O in and density, viscosity of partially CO<sub>2</sub> loaded water-lean amino acid salts:</b>	<a href="https://www.doi.org/10.1016/j.jct.2018.06.021">https://www.doi.org/10.1016/j.jct.2018.06.021</a> <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a> <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C10024972&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C10024972&amp;Units=SI</a>
<b>NIST Webbook:</b>	
<b>Physical Solubility and Diffusivity of N<sub>2</sub>O and CO<sub>2</sub> into Aqueous Solutions:</b>	<a href="https://www.doi.org/10.1021/je0301951">https://www.doi.org/10.1021/je0301951</a>
<b>Diffusion Coefficients of N<sub>2</sub>O in 1,4-Butanediamine Solutions Using a New Method Apparatus for gas solubility measurements and (0.3 to 1.4) mol dm<sup>-3</sup>:</b>	<a href="https://www.doi.org/10.1016/j.jct.2007.05.013">https://www.doi.org/10.1016/j.jct.2007.05.013</a> <a href="https://www.doi.org/10.1016/j.jct.2006.03.013">https://www.doi.org/10.1016/j.jct.2006.03.013</a>
<b>Vapour pressure and excess Gibbs free energy of binary mixtures of hydrogen sulfide with N<sub>2</sub>O and Density and Viscosity of Aqueous Solutions of 1,4-Butanediamine, 2-(Diethylamino)-ethanol, and Their Mixtures from (298.15 to 333.15) K:</b>	<a href="https://www.doi.org/10.1021/je301371p">https://www.doi.org/10.1021/je301371p</a>

## Legend

<b>af:</b>	Acentric Factor
<b>affp:</b>	Proton affinity
<b>basg:</b>	Gas basicity
<b>dm:</b>	Dipole Moment
<b>ea:</b>	Electron affinity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfust:</b>	Enthalpy of fusion at a given temperature
<b>hsubt:</b>	Enthalpy of sublimation at a given temperature
<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>mccvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>pt:</b>	Triple Point Pressure
<b>pvap:</b>	Vapor pressure
<b>rhoc:</b>	Critical density
<b>rhof:</b>	Liquid Density
<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>tt:</b>	Triple Point Temperature
<b>vc:</b>	Critical Volume
<b>zc:</b>	Critical Compressibility
<b>zra:</b>	Rackett Parameter

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