

Ammonia

Other names:	AMMONIA GAS Ammonia, anhydrous Anhydrous ammonia Aromatic Ammonia, Vaporole NH3 NITRO-SIL Spirit of Hartshorn
Inchi:	InChI=1S/H3N/h1H3
InchiKey:	QGZKDVFQNNGYKY-UHFFFAOYSA-N
Formula:	H3N
SMILES:	N
Mol. weight [g/mol]:	17.03
CAS:	7664-41-7

Physical Properties

Property code	Value	Unit	Source
af	0.2500		KDB
affp	853.60	kJ/mol	NIST Webbook
aigt	924.26	K	KDB
basg	819.00	kJ/mol	NIST Webbook
dm	1.50	debye	KDB
fil	15.50	% in Air	KDB
flu	27.00	% in Air	KDB
gf	-16.16	kJ/mol	KDB
gyrad	0.8530		KDB
hf	-45.94 ± 0.35	kJ/mol	NIST Webbook
hf	-45.72	kJ/mol	KDB
ie	10.15	eV	NIST Webbook
ie	10.15 ± 0.01	eV	NIST Webbook
ie	10.82	eV	NIST Webbook
ie	15.02	eV	NIST Webbook
ie	10.16	eV	NIST Webbook
ie	10.16 ± 0.01	eV	NIST Webbook
ie	10.17	eV	NIST Webbook
ie	10.16 ± 0.02	eV	NIST Webbook
ie	14.92	eV	NIST Webbook
ie	10.14	eV	NIST Webbook

ie	10.80	eV	NIST Webbook
ie	10.85	eV	NIST Webbook
ie	11.30	eV	NIST Webbook
ie	10.19	eV	NIST Webbook
ie	10.85	eV	NIST Webbook
ie	14.94 ± 0.03	eV	NIST Webbook
ie	10.18 ± 0.01	eV	NIST Webbook
ie	10.07 ± 0.02	eV	NIST Webbook
ie	10.02	eV	NIST Webbook
ie	10.07 ± 0.01	eV	NIST Webbook
ie	10.20	eV	NIST Webbook
ie	10.07 ± 0.00	eV	NIST Webbook
ie	10.07 ± 0.01	eV	NIST Webbook
ie	10.30 ± 0.00	eV	NIST Webbook
ie	10.45	eV	NIST Webbook
ie	10.20	eV	NIST Webbook
ie	10.18 ± 0.09	eV	NIST Webbook
ie	10.17	eV	NIST Webbook
log10ws	7.60		Crippen Method
logp	0.162		Crippen Method
mcvol	20.840	ml/mol	McGowan Method
nfpaf	%!d(float64=1)		KDB
nfpah	%!d(float64=3)		KDB
pc	11350.00	kPa	KDB
pc	11300.00 ± 5.00	kPa	NIST Webbook
pt	6.06 ± 0.01	kPa	NIST Webbook
pt	6.08	kPa	KDB
rinpol	131.00		NIST Webbook
rinpol	118.00		NIST Webbook
sgb	192.77 ± 0.05	J/mol×K	NIST Webbook
tb	239.82	K	KDB
tc	405.40 ± 0.10	K	NIST Webbook
tc	405.56	K	Critical parameters for ammonia
tc	405.50	K	KDB
tc	405.60 ± 5.00	K	NIST Webbook
tf	195.41	K	KDB
tf	194.95 ± 0.30	K	NIST Webbook
tt	195.41	K	KDB
vc	0.072	m3/kmol	KDB
zc	0.2423820		KDB
zra	0.25		KDB

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpl	89.36	J/mol×K	334.09	Densities, Viscosities, Heat Capacities, and Vapor Liquid Equilibria of Ammonia + Sodium Thiocyanate Solutions at Several Temperatures
cpl	85.83	J/mol×K	324.11	Densities, Viscosities, Heat Capacities, and Vapor Liquid Equilibria of Ammonia + Sodium Thiocyanate Solutions at Several Temperatures
cpl	83.45	J/mol×K	314.14	Densities, Viscosities, Heat Capacities, and Vapor Liquid Equilibria of Ammonia + Sodium Thiocyanate Solutions at Several Temperatures
cpl	82.46	J/mol×K	304.17	Densities, Viscosities, Heat Capacities, and Vapor Liquid Equilibria of Ammonia + Sodium Thiocyanate Solutions at Several Temperatures
hsubt	31.20	kJ/mol	186.00	NIST Webbook
hvapt	22.70	kJ/mol	342.50	NIST Webbook
hvapt	23.50	kJ/mol	220.00	NIST Webbook
hvapt	23.40	kJ/mol	239.00	NIST Webbook

pvap	151.70	kPa	248.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K
pvap	354.80	kPa	268.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K
pvap	429.30	kPa	273.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K
pvap	515.50	kPa	278.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K
pvap	615.10	kPa	283.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K

pvap	728.50	kPa	288.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K
pvap	119.70	kPa	243.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K
pvap	1166.00	kPa	303.12	Vapor Pressures and Liquid Densities of Ammonium Chloride + Ammonia Mixtures
pvap	1553.00	kPa	313.14	Vapor Pressures and Liquid Densities of Ammonium Chloride + Ammonia Mixtures
pvap	2030.00	kPa	323.16	Vapor Pressures and Liquid Densities of Ammonium Chloride + Ammonia Mixtures
pvap	2609.00	kPa	333.13	Vapor Pressures and Liquid Densities of Ammonium Chloride + Ammonia Mixtures
pvap	3305.00	kPa	343.13	Vapor Pressures and Liquid Densities of Ammonium Chloride + Ammonia Mixtures

pvap	4132.00	kPa	353.13	Vapor Pressures and Liquid Densities of Ammonium Chloride + Ammonia Mixtures
pvap	5105.00	kPa	363.12	Vapor Pressures and Liquid Densities of Ammonium Chloride + Ammonia Mixtures
pvap	6243.00	kPa	373.16	Vapor Pressures and Liquid Densities of Ammonium Chloride + Ammonia Mixtures
pvap	290.70	kPa	263.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K
pvap	236.40	kPa	258.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K
pvap	190.20	kPa	253.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K

pvap	857.10	kPa	293.15	The investigation on the vapour liquid phase equilibrium of (ammonia + 1,1,1,2-tetrafluoroethane) system over the temperatures ranging from (243.150 to 283.150) K
rhof	639.00	kg/m3	273.00	
				KDB

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.54557e+01
Coeff. B	-2.33432e+03
Coeff. C	-2.43200e+01
Temperature range (K), min.	195.41
Temperature range (K), max.	405.65

Datasets

Molar heat capacity at constant pressure, J/K/mol

Pressure, kPa - Liquid	Temperature, K - Liquid	Molar heat capacity at constant pressure, J/K/mol - Liquid
1000.00	280.00	76.81
5000.00	280.00	75.62
5000.00	300.00	76.81
5000.00	320.00	77.83
5000.00	360.00	95.20
10000.00	360.00	92.82
15000.00	360.00	90.26

Reference

<https://www.doi.org/10.1016/j.jct.2007.06.013>

Kinematic viscosity, m2/s

Temperature, K - Gas	Pressure, kPa - Gas	Kinematic viscosity, m2/s - Gas
300.00	223.53	0.0000065
300.00	259.87	0.0000055
300.00	329.51	0.0000043
300.00	361.99	0.0000040
300.00	447.86	0.0000031
300.00	465.77	0.0000030
300.00	568.68	0.0000024
300.00	569.79	0.0000025
300.00	676.75	0.0000020
300.00	692.80	0.0000020
300.00	802.37	0.0000017
300.00	802.82	0.0000017
315.00	250.83	0.0000064
315.00	378.83	0.0000042
315.00	506.43	0.0000031
315.00	661.04	0.0000023
315.00	791.02	0.0000019
315.00	938.74	0.0000016
315.00	1080.91	0.0000014
315.00	1284.99	0.0000011
330.00	308.27	0.0000058
330.00	473.84	0.0000037
330.00	664.29	0.0000026
330.00	822.59	0.0000021
330.00	977.41	0.0000017
330.00	1142.12	0.0000014
330.00	1294.82	0.0000013
330.00	1480.49	0.0000011
330.00	1640.80	0.0000010
330.00	1792.05	0.0000009
330.00	1945.84	0.0000008
345.00	297.74	0.0000065
345.00	504.54	0.0000038
345.00	708.88	0.0000027
345.00	914.14	0.0000020
345.00	1123.20	0.0000016
345.00	1332.13	0.0000014

345.00	1537.04	0.0000012
345.00	1744.59	0.0000010
345.00	1955.14	0.0000009
345.00	2315.57	0.0000007
345.00	2530.02	0.0000006
345.00	2769.00	0.0000006
360.00	242.70	0.0000087
360.00	453.28	0.0000047
360.00	661.35	0.0000032
360.00	865.15	0.0000024
360.00	1069.66	0.0000019
360.00	1277.79	0.0000016
360.00	1481.15	0.0000013
360.00	1690.65	0.0000012
360.00	1894.78	0.0000010
360.00	2277.38	0.0000008
360.00	2480.70	0.0000007
360.00	2738.78	0.0000007
360.00	3050.47	0.0000006
360.00	3365.40	0.0000005
375.00	260.57	0.0000101
375.00	483.31	0.0000049
375.00	684.74	0.0000034
375.00	897.66	0.0000026
375.00	1130.68	0.0000020
375.00	1341.71	0.0000016
375.00	1583.75	0.0000014
375.00	1793.43	0.0000012
375.00	2026.26	0.0000010
375.00	2290.42	0.0000009
375.00	2599.92	0.0000008
375.00	2933.13	0.0000007
375.00	3140.48	0.0000006
375.00	3379.20	0.0000006

Reference

<https://www.doi.org/10.1016/j.jct.2007.07.002>

Mass density, kg/m3

Temperature, K - Liquid	Pressure, kPa - Liquid	Mass density, kg/m3 - Liquid
450.00	30000.00	355.3

450.00	40000.00	403.7
450.00	50000.00	433.4
450.00	60000.00	455.2
450.00	70000.00	472.6
450.00	80000.00	487.4
450.00	90000.00	500.2
450.00	100000.00	511.7
450.00	120000.00	531.4
450.00	140000.00	548.2
450.00	160000.00	563.1
450.00	180000.00	576.3
450.00	200000.00	588.2
500.00	50000.00	350.9
500.00	60000.00	384.9
500.00	70000.00	410.1
500.00	80000.00	430.2
500.00	90000.00	446.9
500.00	100000.00	461.5
500.00	120000.00	485.8
500.00	140000.00	505.9
500.00	160000.00	523.1
500.00	180000.00	538.3
500.00	200000.00	551.9

Reference

<https://www.doi.org/10.1016/j.jct.2008.06.009>

Speed of sound, m/s

Temperature, K - Fluid (supercritical or subcritical phases)	Pressure, kPa - Fluid (supercritical or subcritical phases)	Speed of sound, m/s - Fluid (supercritical or subcritical phases)
297.00	620000.00	2741.0
297.00	700000.00	2836.0
297.00	780000.00	2942.0
297.00	880000.00	3023.0
297.00	970000.00	3113.0
297.00	780000.00	2932.0
298.00	370000.00	2376.0
298.00	740000.00	2903.0
474.00	1430000.00	3280.0
476.00	1550000.00	3375.0
476.00	2200000.00	3788.0

477.00	1090000.00	2954.0
479.00	2220000.00	3797.0
678.00	2920000.00	4000.0
678.00	3080000.00	4072.0
678.00	3140000.00	4103.0
679.00	1480000.00	3084.0
679.00	2440000.00	3740.0
679.00	3750000.00	4342.0
680.00	1350000.00	2982.0
680.00	1350000.00	2970.0

Reference

<https://www.doi.org/10.1021/je800312g>

Sources

An Experimental Study of the pVTx Properties for Aqueous Solutions of Ammonia + Ethanol in the Maximum Pressure Region (water + methanol): Investigation on vapor liquid equilibria for binary systems of metal critical parameters for ammonia: [bmim]Zn2Cl5/NH3 by experiment and modified UNI-FAC model: Solubilities of ammonia in basic imidazolium ionic liquids: Vapor Pressures and Liquid Densities of Ammonium Bromide + Ammonia Mixtures. Method:

Effects of association with impurities in ammonia purification: The experimental investigation of the vapour liquid phase equilibrium for NH3 Solubilities and Physical Properties of Ethylamine Hydrochloride Pressure and Temperature Effects on the Ammonia + Water + NaOH and Ammonia + Water + NaOH Solutions: (1 x)H2O} at T = (450 and 500) K over the vapour liquid equilibrium of 200 ammonia + Lithium Nitrate + Water and Ammonia + Measurements of the critical point properties for {xNH3 + (1 x)H2O} with K_D (0.9098, 0.7757, 0.6808):

Densities, Viscosities, Heat Capacities, and Vapor Liquid Equilibria of Ammonia + Sodium Fluoride of ammonia + large liquid mixtures: Suitable alternative working fluids for absorption cycle refrigeration (298.2-353.15 K and 0-300) Squid in Fluid Ammonia to 3.8 GPa and 680 K: The investigation on the vapour liquid phase equilibrium of (ammonia + 2,2,2-trifluoroethanol) system over properties of NH3 in ethylom (243.150-353.15) K: Deep eutectic solvents: Pressure: Solubility of ammonium chloride in a MgCl2-NH4Cl-NH3-H2O system at 298 K: New experiments measuring the solubility of anionic salts in liquid ammonia: Densities, Viscosities, and Heat Capacities of Ammonia + Lithium Nitrate and Ammonia + Lithium Nitrate Solutions of Ammonia in Methanol: (293.15 and 353.15) K:

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Solubilities and Thermodynamic Properties of NH₃ in Glycerin and its Derivatives:
Vapor-Liquid Equilibria of Ammonia + Water + Potassium Hydroxide and Evaluation of Methanesulfonamide-Based Binary Heat Capacities from 293.15 to 333.15 K:
Densities and Liquid Densities of Ammonium Chloride + Ammonia Mixtures:
Liquid Phase Diagram for the System MgCl₂-NH₃-CH₃OH at 298 K: Densities and viscosities of, and NH₃ solubilities in deep eutectic solvents:
Vapor Pressure and Liquid Density of Ammonium Iodide + Ammonia:
Vapor-liquid equilibria of NH₃ in (NH₃ + H₂O) and (NH₃ + PZ + H₂O) system:
Vapor Liquid Equilibrium Measurements of NH₃ + H₂O + Ionic Liquids (DMF, [Dmim]BF₄, and [Dmim]DMP) Systems:
Crippen Method:

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Vapor liquid equilibrium of CO₂ in NH₃ CO₂ SO₂ H₂O system:
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Behavior of the Isochoric Heat Capacity of Ammonia and Density and Viscosity of Aqueous Solutions of Piperazine, Ammonia, and Propylene Glycol:
Properties of CO₂ Absorbents Blended with Ammonia, Ammonia, and Ethanol between 273.15 and 333.15 K:
and 353.15 K:

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Legend

af:	Acentric Factor
affp:	Proton affinity
aigt:	Autoignition Temperature
basg:	Gas basicity
cpl:	Liquid phase heat capacity
dm:	Dipole Moment
fl:	Lower Flammability Limit
flu:	Upper Flammability Limit
gf:	Standard Gibbs free energy of formation
gyrad:	Radius of Gyration
hf:	Enthalpy of formation at standard conditions
hsubt:	Enthalpy of sublimation at a given temperature
hvapt:	Enthalpy of vaporization at a given temperature
ie:	Ionization energy
kvisc:	Kinematic viscosity
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume

nfpaf:	NFPA Fire Rating
nfpah:	NFPA Health Rating
pc:	Critical Pressure
pt:	Triple Point Pressure
pvap:	Vapor pressure
rho:	Liquid Density
rinpol:	Non-polar retention indices
sgb:	Molar entropy at standard conditions (1 bar)
speedsl:	Speed of sound in fluid
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
tt:	Triple Point Temperature
vc:	Critical Volume
zc:	Critical Compressibility
zra:	Rackett Parameter

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