

# sodium hydroxide

Inchi:	InChI=1S/Na.H2O/h;1H2/q+1;/p-1
InchiKey:	HEMHJVSKTPXQMS-UHFFFAOYSA-M
Formula:	HNaO
SMILES:	O[Na]
Mol. weight [g/mol]:	40.00
CAS:	1310-73-2

## Physical Properties

Property code	Value	Unit	Source
affp	1071.80	kJ/mol	NIST Webbook
basg	1044.80	kJ/mol	NIST Webbook

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.51741e+01
Coeff. B	-1.75145e+04
Coeff. C	7.01000e+00
Temperature range (K), min.	773.15
Temperature range (K), max.	1873.15

## Sources

- Solubilities of betulin and betulinic acid in sodium hydroxide aqueous solutions of various molalities at temperatures from 273.2 K to 323.2 K: <https://www.doi.org/10.1016/j.jct.2013.07.014>
- Experimental solid-liquid liquidrite, and sodium calcium oxide NaOH + CaO systems: <https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>
- Measurement and Correlation for the Solubility of Sodium <https://www.doi.org/10.1021/acs.jced.7b01089>
- Experimental solid-liquid liquidrite, and sodium calcium oxide NaOH + CaO systems: <https://www.doi.org/10.1016/j.fluid.2014.07.019>
- Apparent molar heat capacities and apparent molar volumes of zwitterionic, protonated cationic, and deprotonated anionic forms at molalities from (0.002 to 1.0) mol kg<sup>-1</sup>: <https://www.doi.org/10.1016/j.jct.2006.08.008>
- Solubility of betulinic acid and L-glutamic acid at various molalities (0.005 to 0.1 mol kg<sup>-1</sup>) in NaOH and H<sub>2</sub>O systems at 0.1 MPa: <https://www.doi.org/10.1021/acs.jced.8b00400>
- Apparent molar heat capacities and apparent molar volumes of zwitterionic, protonated cationic, and deprotonated anionic forms at molalities from (0.002 to 1.0) mol kg<sup>-1</sup>: <https://www.doi.org/10.1016/j.jct.2006.08.008>

Vapor-Liquid Equilibria of Ammonia + Water + Potassium Hydroxide and Phase Equilibria of the Sodium Hydroxide + Water System at 15 °C and 0.1 MPa. *J. Chem. Eng. Data* 2015, 60, 1031-1034. <https://www.doi.org/10.1021/je049708+>

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Thermodynamic Properties of (Na, K)BO<sub>2</sub> + (Na, K)OH + H<sub>2</sub>O. *J. Chem. Eng. Data* 2015, 60, 1043-1046. <https://www.doi.org/10.1016/j.jct.2005.07.019>

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Thermodynamic Properties of (Na, K)BO<sub>2</sub> + (Na, K)OH + H<sub>2</sub>O. *J. Chem. Eng. Data* 2015, 60, 1059-1062. <https://www.doi.org/10.1021/acs.jced.7b00690>

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Thermodynamic Properties of (Na, K)BO<sub>2</sub> + (Na, K)OH + H<sub>2</sub>O. *J. Chem. Eng. Data* 2015, 60, 1075-1078. <https://www.doi.org/10.1016/j.jct.2011.03.002>

Thermodynamic Properties of (Na, K)BO<sub>2</sub> + (Na, K)OH + H<sub>2</sub>O. *J. Chem. Eng. Data* 2015, 60, 1079-1082. <https://www.doi.org/10.1016/j.jct.2006.05.013>

Thermodynamic Properties of (Na, K)BO<sub>2</sub> + (Na, K)OH + H<sub>2</sub>O. *J. Chem. Eng. Data* 2015, 60, 1083-1086. <https://www.doi.org/10.1016/j.tca.2005.11.035>

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Thermodynamic Properties of (Na, K)BO<sub>2</sub> + (Na, K)OH + H<sub>2</sub>O. *J. Chem. Eng. Data* 2015, 60, 1099-1102. <https://www.doi.org/10.1016/j.jct.2018.08.037>

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## Legend

- affp: Proton affinity
- basg: Gas basicity
- vpap: Vapor pressure

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