

sodium fluoride

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

CAS:

InChI=1S/FH.Na/h1H;/q;+1/p-1

PUZPDOWCWNUUKD-UHFFFAOYSA-M

FNa

F[Na]

41.99

7681-49-4

Physical Properties

Property code	Value	Unit	Source
ea	0.52 ± 0.01	eV	NIST Webbook
ea	1.10 ± 0.20	eV	NIST Webbook
ea	0.42	eV	NIST Webbook
ea	1.12 ± 0.21	eV	NIST Webbook
ea	1.35	eV	NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cps	62.25	J/mol×K	1102.90	Calorimetric analysis of NaF and NaLaF4
cps	58.08	J/mol×K	630.60	Calorimetric analysis of NaF and NaLaF4
cps	58.11	J/mol×K	645.60	Calorimetric analysis of NaF and NaLaF4
cps	59.55	J/mol×K	675.70	Calorimetric analysis of NaF and NaLaF4
cps	59.36	J/mol×K	690.60	Calorimetric analysis of NaF and NaLaF4
cps	58.15	J/mol×K	720.60	Calorimetric analysis of NaF and NaLaF4
cps	59.11	J/mol×K	735.50	Calorimetric analysis of NaF and NaLaF4

cps	58.85	J/mol×K	750.60	Calorimetric analysis of NaF and NaLaF4
cps	59.28	J/mol×K	765.60	Calorimetric analysis of NaF and NaLaF4
cps	58.36	J/mol×K	795.60	Calorimetric analysis of NaF and NaLaF4
cps	58.26	J/mol×K	825.50	Calorimetric analysis of NaF and NaLaF4
cps	58.49	J/mol×K	840.40	Calorimetric analysis of NaF and NaLaF4
cps	59.14	J/mol×K	855.60	Calorimetric analysis of NaF and NaLaF4
cps	58.37	J/mol×K	870.60	Calorimetric analysis of NaF and NaLaF4
cps	59.37	J/mol×K	885.50	Calorimetric analysis of NaF and NaLaF4
cps	59.70	J/mol×K	900.50	Calorimetric analysis of NaF and NaLaF4
cps	60.11	J/mol×K	915.50	Calorimetric analysis of NaF and NaLaF4
cps	59.64	J/mol×K	945.50	Calorimetric analysis of NaF and NaLaF4
cps	60.43	J/mol×K	960.60	Calorimetric analysis of NaF and NaLaF4
cps	59.89	J/mol×K	975.40	Calorimetric analysis of NaF and NaLaF4
cps	61.01	J/mol×K	990.50	Calorimetric analysis of NaF and NaLaF4
cps	59.76	J/mol×K	907.50	Calorimetric analysis of NaF and NaLaF4
cps	60.21	J/mol×K	922.30	Calorimetric analysis of NaF and NaLaF4
cps	60.58	J/mol×K	937.40	Calorimetric analysis of NaF and NaLaF4
cps	59.98	J/mol×K	952.50	Calorimetric analysis of NaF and NaLaF4
cps	60.92	J/mol×K	967.50	Calorimetric analysis of NaF and NaLaF4

cps	60.11	J/mol×K	982.60	Calorimetric analysis of NaF and NaLaF4
cps	60.91	J/mol×K	997.50	Calorimetric analysis of NaF and NaLaF4
cps	61.02	J/mol×K	1012.60	Calorimetric analysis of NaF and NaLaF4
cps	60.56	J/mol×K	1027.60	Calorimetric analysis of NaF and NaLaF4
cps	61.86	J/mol×K	1042.60	Calorimetric analysis of NaF and NaLaF4
cps	61.63	J/mol×K	1057.70	Calorimetric analysis of NaF and NaLaF4
cps	62.08	J/mol×K	1072.70	Calorimetric analysis of NaF and NaLaF4
cps	62.41	J/mol×K	1087.80	Calorimetric analysis of NaF and NaLaF4
srf	0.19	N/m	1325.00	Surface tension of light rare earth fluoride molten salts electrolytesystem

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.78327e+01
Coeff. B	-2.62688e+04
Coeff. C	5.17000e+00
Temperature range (K), min.	1269.00
Temperature range (K), max.	2060.00

Sources

Activity coefficients of NaF in aqueous mixtures with e-increasing co-solvent: NaOH and NaOH + NaF and NaOH + NaF + NaF at 298.15 K: <https://www.doi.org/10.1016/j.fluid.2016.05.019>
 Calorimetric analysis of NaF and NaLaF4: <https://www.doi.org/10.1016/j.jct.2006.03.002>

Enthalpies of Dilution of
(2S,3R,4R,5R)-Hexane-1,2,3,4,5,6-hexol
Surface Tension and Density in Hande
NaF-Aqueous Based Electrolyte (18.15) K:
Surface tension of light rare earth
fluoride molten salts electrolytesystem:
Solubility measurements in
Na-F-CO₃-HCO₃-H₂O system at (308.15
and 323.15) K and development of a
Pressurized equilibrium model for the
Na-F-CO₃-SO₄-CO₃-HCO₃-H₂O system:
Partial molar volume and partial molar
compressibility of four homologous
monovalent acids in aqueous sodium
fluoride solutions at different
temperatures
Physicochemical properties of the
system(LiF + NaF + KF(eut.) +
Electrolyte Conductivity of Electrolytes
Found in Natural Waters, isom (5, 10, 20)
dependence of the
Density of Aqueous Alkali Halide Salt
Solutions by Experiment and Molecular
simulation: sodiumhydroxide + water
Systems: experimental and calculation:
+ water) and (sucrose + water) mixtures
at 298.15 K:

<https://www.doi.org/10.1021/je1007394>
<https://www.doi.org/10.1021/je2005825>
<https://www.doi.org/10.1016/j.tca.2016.03.033>
<https://www.doi.org/10.1016/j.jct.2018.10.029>
<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>
<https://www.doi.org/10.1016/j.jct.2011.01.004>
<http://webbook.nist.gov/cgi/cbook.cgi?ID=C7681494&Units=SI>
<https://www.doi.org/10.1016/j.jct.2014.03.024>
<https://www.doi.org/10.1021/je101012n>
<https://www.doi.org/10.1021/je500420g>
<https://www.doi.org/10.1016/j.tca.2013.06.008>
<https://www.doi.org/10.1016/j.jct.2004.07.002>

Legend

cps: Solid phase heat capacity
ea: Electron affinity
pvap: Vapor pressure
srf: Surface Tension

Latest version available from:

<https://www.cheméo.com/cid/13-766-7/sodium-fluoride.pdf>

Generated by Cheméo on 2025-12-05 10:58:55.808309717 +0000 UTC m=+4680533.338350381.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.