

potassium

Inchi:	InChI=1S/K
InchiKey:	ZLMJMSJWJFRBEC-UHFFFAOYSA-N
Formula:	K
SMILES:	[K]
Mol. weight [g/mol]:	39.10
CAS:	7440-09-7

Physical Properties

Property code	Value	Unit	Source
hf	89.00 ± 0.80	kJ/mol	NIST Webbook
ie	4.50 ± 1.00	eV	NIST Webbook
ie	4.34	eV	NIST Webbook
ie	4.10 ± 0.30	eV	NIST Webbook
ie	4.40	eV	NIST Webbook
ie	4.34 ± 0.00	eV	NIST Webbook
ie	4.34	eV	NIST Webbook
ie	4.34	eV	NIST Webbook
ie	4.34	eV	NIST Webbook
ie	4.34 ± 0.00	eV	NIST Webbook
nfpaf	%!d(float64=1)		KDB
nfpah	%!d(float64=3)		KDB
nfpas	%!d(float64=2)		KDB
sgb	160.34 ± 0.00	J/mol×K	NIST Webbook
ss	64.68 ± 0.20	J/mol×K	NIST Webbook
tb	1037.00 ± 1.00	K	NIST Webbook
tt	336.35 ± 0.10	K	NIST Webbook

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{\text{vp}}) = A + B/(T + C)$
Coeff. A	1.40440e+01
Coeff. B	-9.54174e+03

Coeff. C	-1.86800e+01
Temperature range (K), min.	336.35
Temperature range (K), max.	2223.00

Sources

Low-Temperature Heat Capacities and Standard Molar Enthalpy of Formation NIST Webbook
Hydrogen Phthalate C8H5KO4(s):
A calorimetric and thermodynamic investigation of potassium uranyl low-temperature heat capacities and standard molar enthalpy of formation of potassium uranyl oxo-phthalate ($\text{UO}_2\text{C}_8\text{H}_5\text{O}_4\text{K}(s)$): microporous materials for two boron ceramics. <https://doi.org/10.1016/j.jct.2012.09.014>
Investigation of $[(\text{UO}_2)(\text{MoO}_4)\text{O}_2]$ compounds with Astatine Rb and $\text{A}_{2}\text{K}_{2}\text{U}_{2}(\text{WO}_4)_2$ compounds with A system $\text{K}_2\text{UW}_2\text{O}_9$ and $\text{K}_2\text{UW}_2\text{O}_9$ phase relations in the $\text{K}_2\text{WO}_4-\text{UO}_3-\text{H}_2\text{O}$ and $\text{K}_2\text{MoO}_4-\text{K}_2\text{WO}_4-\text{UO}_3-\text{H}_2\text{O}$ systems: Pressure: Synthesis and energetics of Na, K, Rb and Cs salts by reaction with The thermochemistry of Potassium: Strontium Tetraborate Decahydrated:

- <https://www.doi.org/10.1021/je900759f>
- <http://webbook.nist.gov/cgi/cbook.cgi?ID=C7440097&Units=SI>
- <https://www.doi.org/10.1016/j.jct.2012.09.014>
- <https://www.doi.org/10.1007/s10765-009-0568-4>
- <https://www.doi.org/10.1016/j.jct.2015.08.032>
- <https://www.doi.org/10.1016/j.jct.2015.06.028>
- <https://www.doi.org/10.1016/j.jct.2017.03.039>
- <https://www.cheric.org/research/kdb/hcprop/showprop.php?cmpid=1956>
- <https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>
- <https://www.doi.org/10.1016/j.jct.2017.07.025>
- <https://www.doi.org/10.1016/j.tca.2007.08.004>

Legend

hf:	Enthalpy of formation at standard conditions
ie:	Ionization energy
nfpaf:	NFPA Fire Rating
nfpah:	NFPA Health Rating
nfpas:	NFPA Safety Rating
pvap:	Vapor pressure
sgb:	Molar entropy at standard conditions (1 bar)
ss:	Solid phase molar entropy at standard conditions
tb:	Normal Boiling Point Temperature
tt:	Triple Point Temperature

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