

potassium hydroxide

| | |
|----------------------|-------------------------------|
| Inchi: | InChI=1S/K.H2O/h;1H2/q+1;/p-1 |
| InchiKey: | KWYUFKZDYYNOTN-UHFFFAOYSA-M |
| Formula: | HKO |
| SMILES: | O[K] |
| Mol. weight [g/mol]: | 56.11 |
| CAS: | 1310-58-3 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|-------------|--------|--------------|
| affp | 1101.80 | kJ/mol | NIST Webbook |
| basg | 1075.40 | kJ/mol | NIST Webbook |
| ie | 7.50 ± 1.00 | eV | NIST Webbook |
| ie | 7.50 ± 0.15 | eV | NIST Webbook |

Correlations

| Information | Value |
|-----------------------------|--------------------------------------|
| Property code | pvap |
| Equation | $\ln(P_{\text{vp}}) = A + B/(T + C)$ |
| Coeff. A | 1.42925e+01 |
| Coeff. B | -1.45074e+04 |
| Coeff. C | -1.00390e+02 |
| Temperature range (K), min. | 679.00 |
| Temperature range (K), max. | 1600.00 |

Sources

Solubility of Sodium Oxalate in Concentrated Electrolyte Solutions: Densities and Heat Capacities of the Ammonia + Water + NaOH and Thermodynamics of KOH, NaOH and Cs alkylated phenoxides:
NIST Webbook:

<https://www.doi.org/10.1021/acs.jced.7b00690>
<https://www.doi.org/10.1021/je050512z>
<https://www.doi.org/10.1016/j.tca.2005.02.027>
<http://webbook.nist.gov/cgi/cbook.cgi?ID=C1310583&Units=SI>

Measurement and modelling of solubility for calcium sulfate dihydrate thermal stability characteristics of ROH reactions with KH(PO₃H):
Solubility of the System KOH K₂CrO₄ Al₂O₃ H₂O at 150 deg C in a High Alkali Concentration Determination of the KOH + K₂CrO₄ + H₂O System:
Density Calculations for (Na, K)BH₄ + (Na, K)BO₂ + (Na, K)OH + H₂O
Solubility of Calcium Hydrogen Arsenite
Binary Systems Containing Selected Heteroatoms: Property, Landolt and Solubility of aqueous amino acid salt and amine Ammonium Salt Solutions at Vapor Pressure:
Vapor-Liquid Equilibria of Ammonia + Water + Potassium Hydroxide and Ammonia + Water + Sodium Hydroxide Solutions at Temperatures from (293.15 to 353.15) K:

Legend

| | |
|--------------|-------------------|
| affp: | Proton affinity |
| basg: | Gas basicity |
| ie: | Ionization energy |
| pvap: | Vapor pressure |

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- <https://www.doi.org/10.1016/j.fluid.2010.06.012>
<https://www.doi.org/10.1016/j.tca.2012.08.020>
<https://www.doi.org/10.1021/je3004782>
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<https://www.doi.org/10.1021/je200216n>
<https://www.doi.org/10.1021/je700315u>
<https://www.doi.org/10.1016/j.jct.2011.09.012>
<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>
<https://www.doi.org/10.1021/je049708+>