



## ***Corrigendum to*** **“Compilation of Henry’s law constants (version 4.0) for water as solvent” published in Atmos. Chem. Phys., 15, 4399–4981, 2015**

**Rolf Sander**

Atmospheric Chemistry Department, Max Planck Institute for Chemistry, P.O. Box 3060, 55020 Mainz, Germany

**Correspondence:** Rolf Sander (rolf.sander@mpic.de)

Published: 22 February 2021

Since the compilation of Henry’s law constants was published in 2015, a couple of errors have been discovered, as summarized below. If additional errata become necessary in the future, they will be posted on my Henry’s law website at <http://www.henrys-law.org> (last access: 19 February 2021).

- Due to a unit conversion error, all  $H'$  values (products of Henry’s law and acidity constants) for strong acids in my compilation were off by a factor of 1000. The corrected data are shown in Table 1.
- When analyzing Henry’s law constant for  $\text{N}_2\text{O}_5$  from Fried et al. (1994), I had incorrectly assumed that they were referring to Henry’s law constant  $H^{cp}$  (in  $\text{M atm}^{-1}$ ). Re-checking their publication, however, I noticed that they used the dimensionless Henry’s law constant  $H^{cc}$ . Thus, the correct conversion yields  $H^{cp}(\text{N}_2\text{O}_5) = 8.7 \times 10^{-4} \text{ mol (m}^3 \text{ Pa)}^{-1}$  and a temperature dependence of  $d \ln H^{cp} / d(1/T) = 3600 \text{ K}$ .

- I adopted an incorrect value for HOBr from the JPL data evaluation by Sander et al. (2011). The corrected value, based on the updated JPL data evaluation by Burkholder et al. (2015), is  $H^{cp}(\text{HOBr}) > 13 \text{ mol (m}^3 \text{ Pa)}^{-1}$ .
- The reference “Lia et al. (2004)” should be Li et al. (2004).

**Table 1.** Corrected values of  $H' = H^{CP} \times K_A$  for strong acids, where  $H^{CP}$  is Henry's law constant, and  $K_A$  is the acidity constant.

|                      | Corrected value   |   | Reference                     |
|----------------------|---|---|-------------------------------|
| $H'(\text{HNO}_3) =$ | $2.6 \times 10^7 \times \exp\left(8700 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$       | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Clegg and Brimblecombe (1990) |
| $H'(\text{HNO}_3) =$ | $2.4 \times 10^7 \times \exp\left(8700 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$       | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Brimblecombe and Clegg (1989) |
| $H'(\text{HF}) =$    | $9.4 \times 10^1 \times \exp\left(7400 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$       | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Brimblecombe and Clegg (1989) |
| $H'(\text{HCl}) =$   | $2.0 \times 10^7$   | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Clegg and Brimblecombe (1986) |
| $H'(\text{HCl}) =$   | $2.0 \times 10^7 \times \exp\left(9000 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$       | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Wagman et al. (1982)          |
| $H'(\text{HCl}) =$   | $2.0 \times 10^7 \times \exp\left(9000 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$       | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Carslaw et al. (1995)         |
| $H'(\text{HCl}) =$   | $2.0 \times 10^7 \times \exp\left(9000 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$       | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Brimblecombe and Clegg (1989) |
| $H'(\text{HBr}) =$   | $7.0 \times 10^9 \times \exp\left(10\,000 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$    | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Wagman et al. (1982)          |
| $H'(\text{HBr}) =$   | $7.1 \times 10^9 \times \exp\left(6100 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$       | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Chameides and Stelson (1992)  |
| $H'(\text{HBr}) =$   | $8.2 \times 10^9 \times \exp\left(10\,000 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$    | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Carslaw et al. (1995)         |
| $H'(\text{HBr}) =$   | $1.3 \times 10^{10} \times \exp\left(10\,000 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$ | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Brimblecombe and Clegg (1989) |
| $H'(\text{HI}) =$    | $2.1 \times 10^{10} \times \exp\left(9800 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$    | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Wagman et al. (1982)          |
| $H'(\text{HI}) =$    | $2.5 \times 10^{10} \times \exp\left(9800 \text{ K} \left(\frac{1}{T} - \frac{1}{T^\ominus}\right)\right)$    | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Brimblecombe and Clegg (1989) |
| $H'(\text{MSA}) =$   | $6.4 \times 10^{14}$  | $\text{mol}^2 (\text{m}^6 \text{ Pa})^{-1}$ | Brimblecombe and Clegg (1988) |

*Acknowledgements.* I would like to thank Yuling Chen for pointing out an inconsistency in the data for HF, which led to the discovery of the unit conversion error for the  $H'$  values of strong acids.

## References

- Brimblecombe, P. and Clegg, S. L.: The solubility and behaviour of acid gases in the marine aerosol, *J. Atmos. Chem.*, 7, 1–18, <https://doi.org/10.1007/BF00048251>, 1988.
- Brimblecombe, P. and Clegg, S. L.: Erratum, *J. Atmos. Chem.*, 8, 95, <https://doi.org/10.1007/BF00053818>, 1989.
- Burkholder, J. B., Sander, S. P., Abbatt, J., Barker, J. R., Huie, R. E., Kolb, C. E., Kurylo, M. J., Orkin, V. L., Wilmouth, D. M., and Wine, P. H.: Chemical Kinetics and Photochemical Data for Use in Atmospheric Studies, Evaluation No. 18, JPL Publication 15-10, Jet Propulsion Laboratory, Pasadena, available at: <http://jpldataeval.jpl.nasa.gov> (last access: 19 February 2021), 2015.
- Carslaw, K. S., Clegg, S. L., and Brimblecombe, P.: A thermodynamic model of the system  $\text{HCl-HNO}_3\text{-H}_2\text{SO}_4\text{-H}_2\text{O}$ , including solubilities of HBr, from < 200 to 328 K, *J. Phys. Chem.*, 99, 11557–11574, <https://doi.org/10.1021/J100029A039>, 1995.
- Chameides, W. L. and Stelson, A. W.: Aqueous phase chemical processes in deliquescent sea-salt aerosols: A mechanism that couples the atmospheric cycles of S and sea salt, *J. Geophys. Res.*, 97D, 20565–20580, <https://doi.org/10.1029/92JD01923>, 1992.
- Clegg, S. L. and Brimblecombe, P.: The dissociation constant and Henry's law constant of HCl in aqueous solution, *Atmos. Environ.*, 20, 2483–2485, [https://doi.org/10.1016/0004-6981\(86\)90079-X](https://doi.org/10.1016/0004-6981(86)90079-X), 1986.
- Clegg, S. L. and Brimblecombe, P.: Equilibrium partial pressures and mean activity and osmotic coefficients of 0–100 % nitric acid as a function of temperature, *J. Phys. Chem.*, 94, 5369–5380, <https://doi.org/10.1021/J100376A038>, 1990.
- Fried, A., Henry, B. E., Calvert, J. G., and Mozurkewich, M.: The reaction probability of  $\text{N}_2\text{O}_5$  with sulfuric acid aerosols at stratospheric temperatures and compositions, *J. Geophys. Res.*, 99D, 3517–3532, <https://doi.org/10.1029/93JD01907>, 1994.
- Li, S., Chen, Z., and Shi, F.: Determination of Henry's Law constant for methyl hydroperoxide by long path FTIR, *Prog. Nat. Sci.*, 14, 765–769, <https://doi.org/10.1080/10020070412331344291>, 2004.
- Sander, S. P., Abbatt, J., Barker, J. R., Burkholder, J. B., Friedl, R. R., Golden, D. M., Huie, R. E., Kolb, C. E., Kurylo, M. J., Moortgat, G. K., Orkin, V. L., and Wine, P. H.: Chemical Kinetics and Photochemical Data for Use in Atmospheric Studies, Evaluation No. 17, JPL Publication 10-6, Jet Propulsion Laboratory, Pasadena, available at: <http://jpldataeval.jpl.nasa.gov> (last access: 19 February 2021), 2011.
- Wagman, D. D., Evans, W. H., Parker, V. B., Schumm, R. H., Halow, I., Bailey, S. M., Churney, K. L., and Nuttall, R. L.: The NBS tables of chemical thermodynamic properties; Selected values for inorganic and  $\text{C}_1$  and  $\text{C}_2$  organic substances in SI units, *J. Phys. Chem. Ref. Data*, 11, suppl. 2, 1982.