

## Chemical investigation of hassium (Hs, Z=108)

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The heaviest element, whose chemical behavior has been studied so far is bohrium (Bh) with  $Z=107$  [1] behaving like a typical member of group 7 of the periodic table. The longest-lived  $\alpha$ -decaying isotope of the next heavier element hassium (Hs,  $Z=108$ ) is  $^{269}\text{Hs}$  ( $T_{1/2}=11.3$  s) which has been identified in the decay chain of  $^{277}\text{112}$  [2,3]. Hs is supposed to be a member of group 8 of the periodic table and should thus form a very volatile tetroxide. Relativistic density functional calculations predicted the electronic structure of  $\text{HsO}_4$  to be similar to the one of  $\text{OsO}_4$  [4]. Application of different semiempirical models of the interaction of a  $\text{MeO}_4$  molecule with quartz surface predicted the adsorption behavior of  $\text{OsO}_4$  and  $\text{HsO}_4$  to be very similar [4]. Extrapolations of trends within group 8 of the periodic table also predicted  $\text{HsO}_4$  and  $\text{OsO}_4$  to behave similar in a gas adsorption chromatography experiment [5].

Hs isotopes were produced directly in the reaction  $^{248}\text{Cm}(^{26}\text{Mg};5,4n)^{269,270}\text{Hs}$  at the UNILAC at GSI Darmstadt [6]. Hs isotopes recoiling from the target were thermalized and oxidized in a  $\text{He}/\text{O}_2$  mixture in the recoil chamber of the In-situ Volatilization and On-line detection apparatus IVO [7]. Volatile  $\text{HsO}_4$  was transported with the carrier gas to the Cryo-On-Line-Detector (COLD), a thermochromatography device. Along a narrow channel formed of PIN-diodes registering  $\alpha$ -decaying and spontaneously fissioning (SF) nuclides, a temperature gradient from  $-20$  to  $-170$  °C was established. The deposition temperature of volatile species could therefore be determined, allowing for the determination of their adsorption enthalpy. COLD is an improved version of the Cryo-Thermochromatography Separator CTS developed at Berkeley [8].

Five decay chains were detected in the course of the experiment which were attributed to  $^{269}\text{Hs}$  or the so far unknown isotope  $^{270}\text{Hs}$  [6]. In addition, two  $\alpha$ -SF correlations were observed in detectors 3 and 4 that still have a rather low random probability, but could not be assigned with certainty to either  $^{269}\text{Hs}$  or  $^{270}\text{Hs}$  [6]. The deposition temperature of the Hs containing molecules was determined to  $(-44\pm 5)$  °C giving strong evidence of the formation of  $\text{HsO}_4$ . In an irradiation of a  $^{152}\text{Gd}$  target,  $^{172}\text{Os}$  ( $T_{1/2}=19.2$  s) was produced in the reaction  $^{152}\text{Gd}(^{26}\text{Mg};6n)$  and a deposition temperature of  $(-82\pm 5)$  °C was measured for  $^{172}\text{OsO}_4$ . The deposition distribution in the COLD array along the detector pairs is shown in Fig. 1. From these deposition peaks the adsorption enthalpies were deduced applying a Monte-Carlo simulation based on a microscopic description of the transport process in the chromatography column [9], i.e. in the COLD system. Since the half-life of the nuclide is a crucial parameter in this simulation and this value has not yet been measured for  $^{270}\text{Hs}$ , only the three events

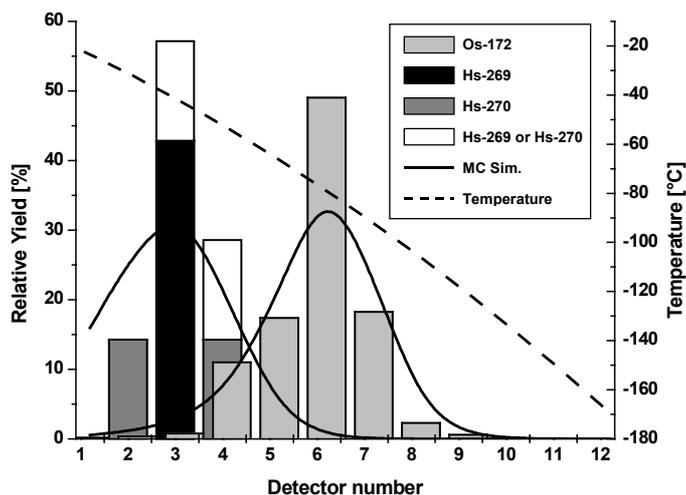


Fig. 1 Merged thermochromatograms of  $\text{OsO}_4$  and  $\text{HsO}_4$ . The solid lines represent results of a Monte-Carlo Simulation with  $\Delta H_{\text{ads}}$  values of  $-39.5$   $\text{kJ}\cdot\text{mol}^{-1}$  ( $\text{OsO}_4$ ) and  $-47$   $\text{kJ}\cdot\text{mol}^{-1}$  ( $\text{HsO}_4$ ), respectively. The dashed line indicates the temperature gradient.

assigned to  $^{269}\text{Hs}$  were used for the simulation.  $\Delta H_{\text{ads}}(\text{HsO}_4)=(-47\pm 2)$   $\text{kJ}\cdot\text{mol}^{-1}$  (68 % c.i.) was evaluated, compared to  $\Delta H_{\text{ads}}(\text{OsO}_4)=(-39.5\pm 1.0)$   $\text{kJ}\cdot\text{mol}^{-1}$ . The latter value is in good agreement with  $\Delta H_{\text{ads}}(\text{OsO}_4)=(-38.0\pm 1.5)$   $\text{kJ}\cdot\text{mol}^{-1}$  evaluated in earlier experiments.

With the formation of a very volatile oxide, presumably  $\text{HsO}_4$ , Hs behaves similar to Os, its next lighter homologue in group 8 of the periodic table.

## References

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