

# Chemistry of flerovium – current status

Lotte Lens<sup>1,2</sup> for the TASCA Flerovium Chemistry Collaboration

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TASCA Workshop 16 – Darmstadt, Germany, August 26, 2016

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# The TASCA Flerovium Chemistry Collaboration

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<b>Univ. Oslo (N)</b>	J.P. Omtvedt
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<b>JAEA, Tokai (J)</b>	M. Asai, Y. Nagame, T. Sato, K. Tsukada
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<b>SINP Kolkata (IN)</b>	S. Lahiri
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<b>ANU Canberra, (AU)</b>	D. Hinde, E. Williams
<b>ITE Warsaw (PL)</b>	M. Wegrzecki

# Acknowledgments

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- ❖ GSI Ion source and accelerator operation staff
- ❖ Electrical / mechanical workshops
- ❖ Experimental electronic department GSI

- ❖ BMBF Verbundforschung



- ❖ JAEA ASRC's Reimei Program
- ❖ ENSAR Travel support Program

# Introduction

The image shows a periodic table of elements. The elements are arranged in rows and columns. The atomic numbers 1 through 118 are shown in the bottom left corner of each element's box. The elements are labeled with their chemical symbols. The elements 112, 114, and 118 are highlighted with red boxes. The element 14 is highlighted in green. The element 118 is labeled as Og (Oganesson).

1																			18
H																			He
Li	Be																		
Na	Mg																		
K	Ca	Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn	Ga	Ge	As	Se	Br	Kr		
Rb	Sr	Y	Zr	Nb	Mo	Tc	Ru	Rh	Pd	Ag	Cd	In	Sn	Sb	Te	I	Xe		
Cs	Ba	La*	Hf	Ta	W	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	At	Rn		
Fr	Ra	Ac <sup>+</sup>	Rf	Db	Sg	Bh	Hs	Mt	Ds	Rg	Cn	(Nh)	Fl	(Mc)	Lv	(Ts)	Og		

## Are elements 112, 114, and 118 relatively inert gases?

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(Received 14 April 1975)*

The Journal of Chemical Physics, Vol. 63, No. 2, 15 July 1975

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# Predictions of adsorption Enthalpies for

Adsorption enthalpies  $-\Delta H_{\text{ads}}$  (in kJ/mol) of Hg-Au<sub>n</sub>, Pb-Au<sub>n</sub>, Cn-Au<sub>n</sub> and Fl-Au<sub>n</sub> for Au(111) and of Hg-SiO<sub>2</sub>, Pb-SiO<sub>2</sub>, Cn-SiO<sub>2</sub>, Fl-SiO<sub>2</sub> for SiO<sub>2</sub> hydroxylated  $\alpha$ -quartz (001) surface

## Hg

	$-\Delta H_{\text{ads}}$ Theory	$-\Delta H_{\text{ads}}$ Experiment	Ref.
SiO <sub>2</sub>	54	42±2	4,2,3
Au	54	98±3	1,7

## Pb

	$-\Delta H_{\text{ads}}$ Theory	$-\Delta H_{\text{ads}}$ Experiment	Ref.
SiO <sub>2</sub>	152	165±4	4,2,3
Au	232	234	1,5,6

## Cn

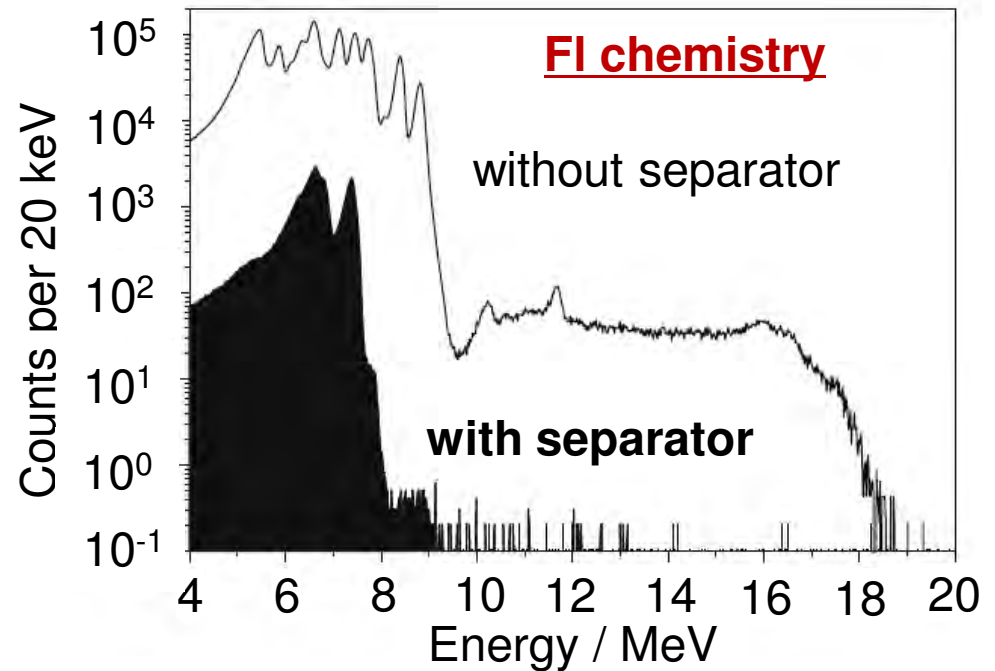
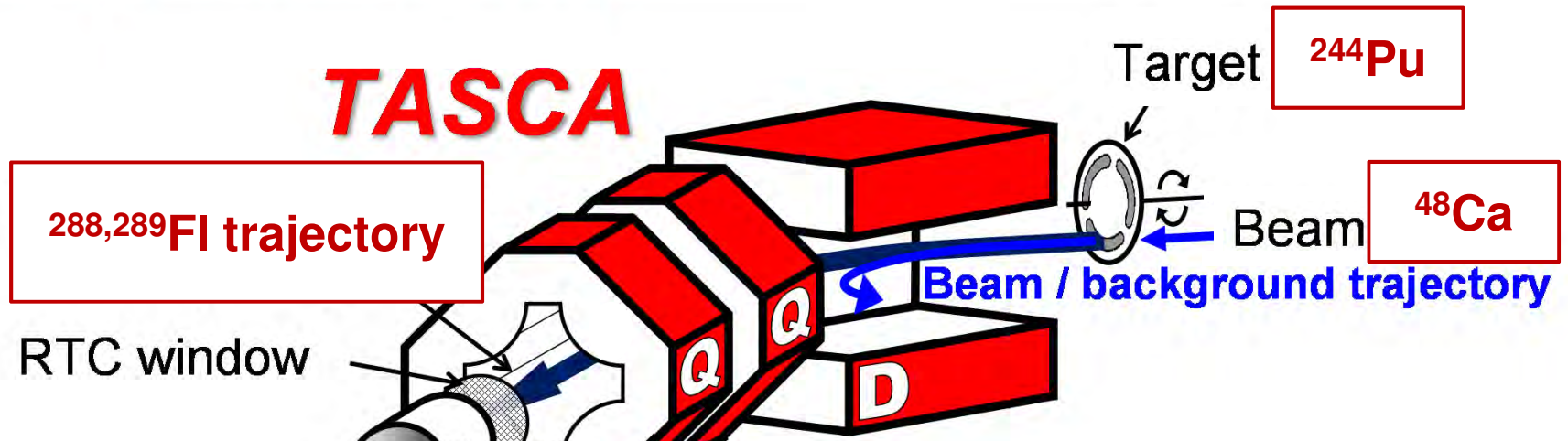
	$-\Delta H_{\text{ads}}$ Theory	$-\Delta H_{\text{ads}}$ Experiment	Ref.
SiO <sub>2</sub>	-38	-	4
Au	45	52 <sup>+4</sup> <sub>-3</sub>	1,5,6

## Fl

	$-\Delta H_{\text{ads}}$ Theory	$-\Delta H_{\text{ads}}$ Experiment	Ref.
SiO <sub>2</sub>	-22	-	4
Au	68	Under discussion	1,5,6

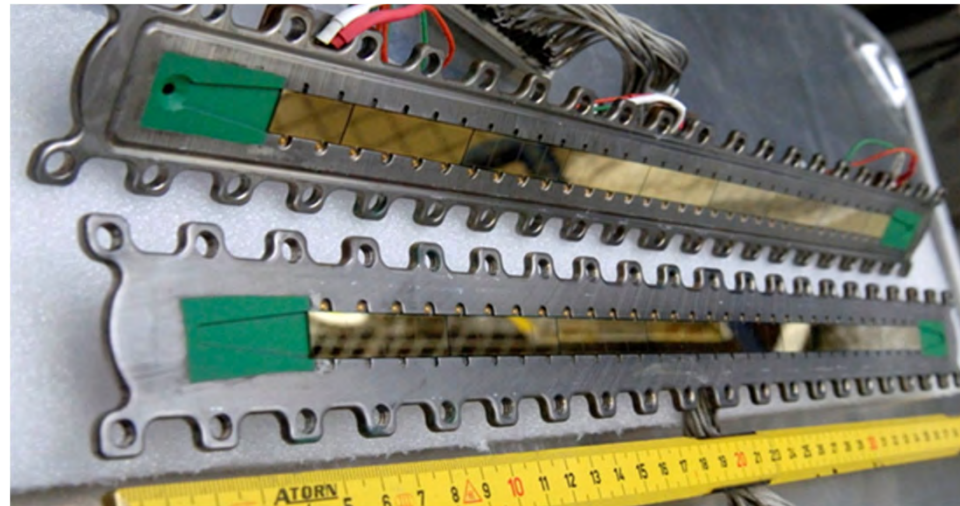
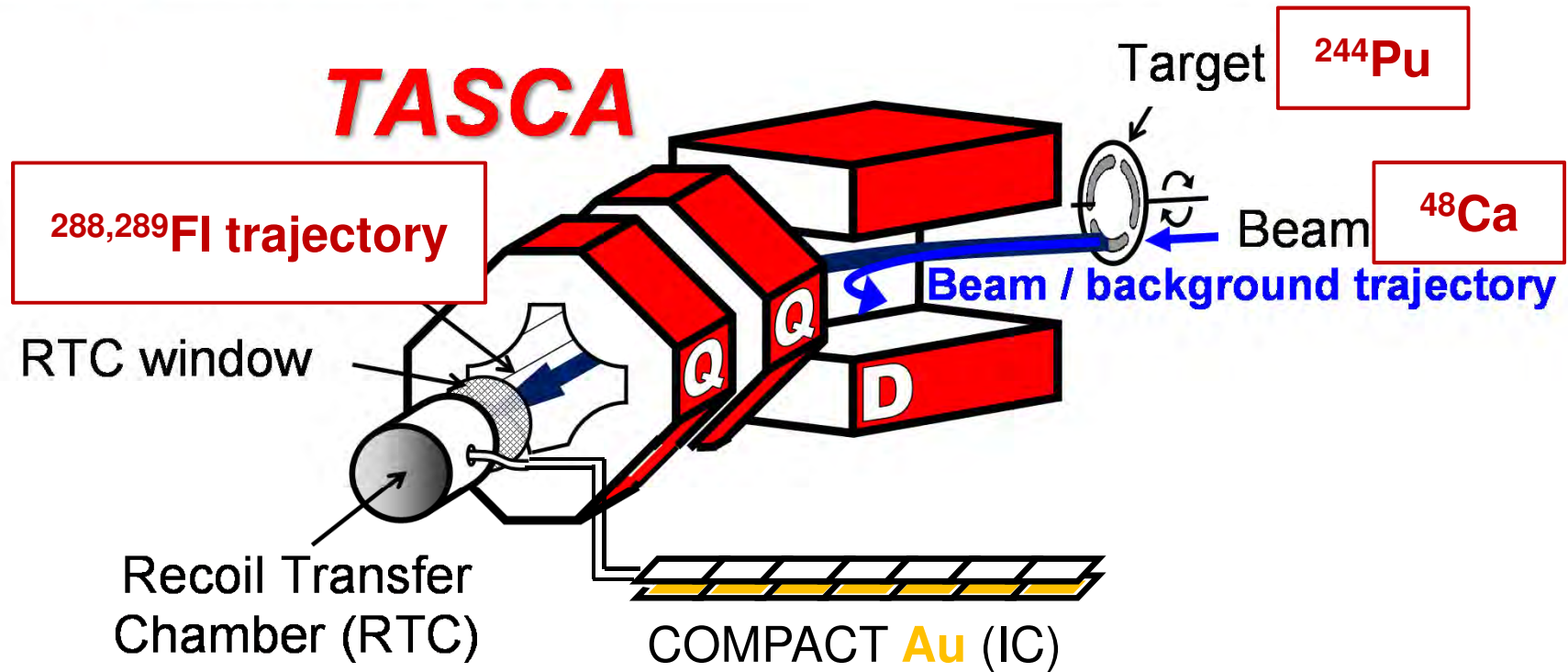
1. V. Pershina *et al.*, J. Chem. Phys. 131, 084713 (2009)
2. S. Soverna, Doctoral Thesis, University Bern (2004)
3. R. Eichler, private communication, (2015) to be published
4. V. Pershina, Phys. Chem. Chem. Phys. 18, 17750-17756 (2016)
5. R. Eichler *et al.*, Nature 447, 72 (2007)
6. R. Eichler *et al.*, Radiochim. Acta 98, 133 (2010)
7. S. Soverna *et al.*, RCA 93 (2005)

# Experimental setup

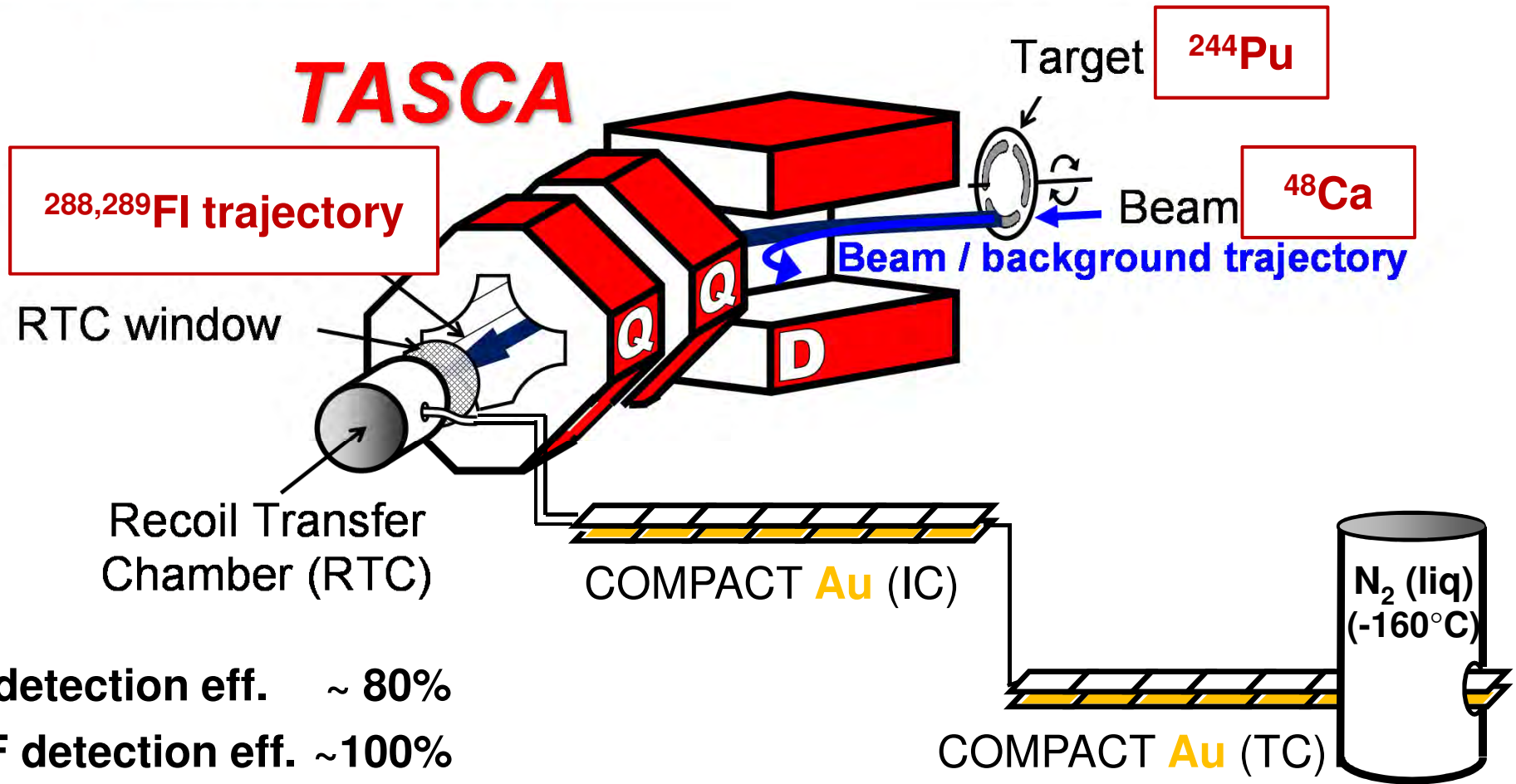




# Experimental setup



# Experimental setup



- ❖  $\alpha$  detection eff.  $\sim 80\%$
- ❖ SF detection eff.  $\sim 100\%$
- ❖ Full digital electronics

[N. Kurz *et al.*, GSI Scientific Report 252 (2012)]

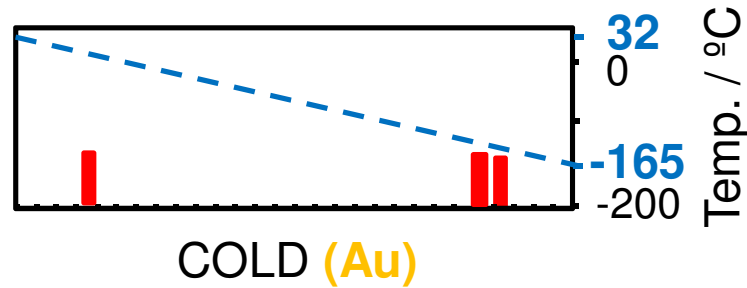
[J. Hoffmann *et al.*, GSI Scientific Report 253 (2012)]

[J. Khuyagbaatar *et al.*, GSI Scientific Report 212 (2012)]



# Experimental status

**2007 (PSI / FLNR)** (without separator, 5 weeks, beam dose:  $6 \cdot 10^{18}$ )

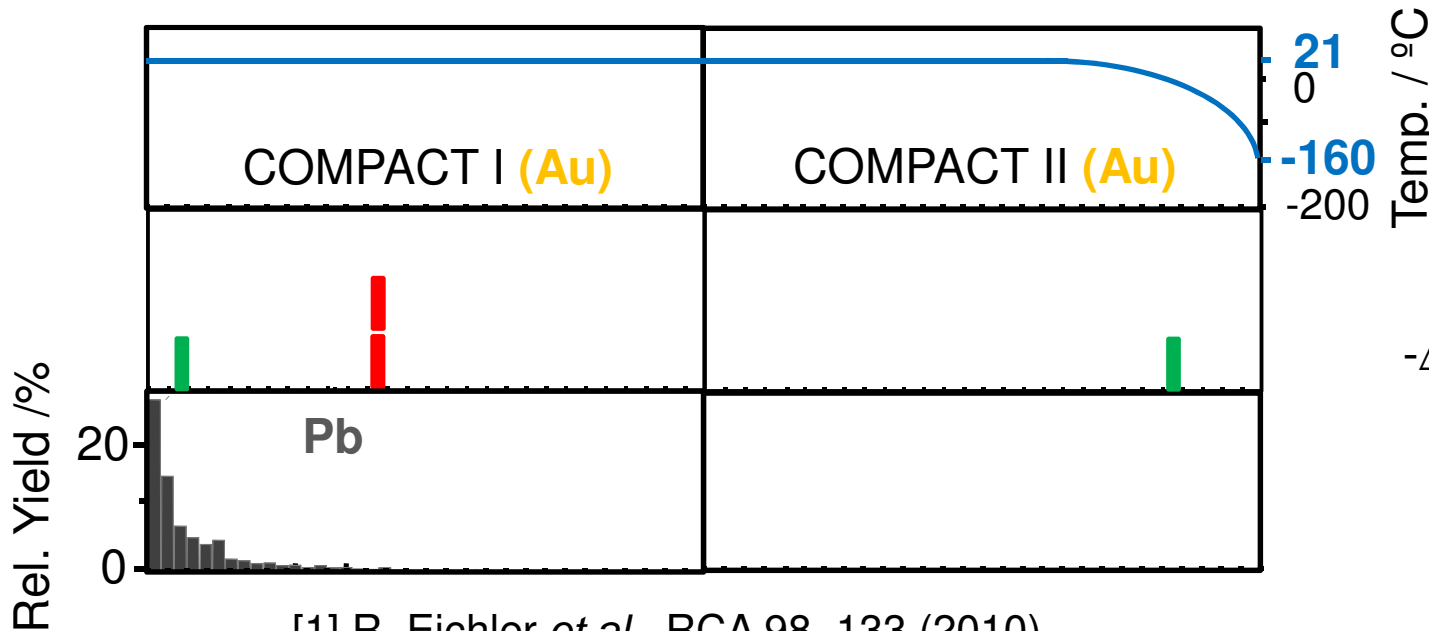


$$-\Delta H_{\text{ads}}^{\text{Au}}(\text{FI}) = 34^{+54}_{-11} \text{ kJ/mol}^{[1]}$$

➤ Weak physisorption bond FI(Au)

**2009 GSI Darmstadt** (with separator, 4 weeks, beam dose:  $4 \cdot 10^{18}$ )

FI decay SF



$$-\Delta H_{\text{ads}}^{\text{Au}}(\text{FI}) > 48 \text{ kJ/mol}^{[2]}$$

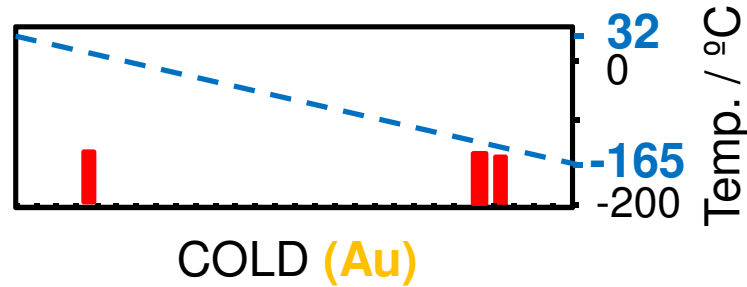
➤ Metallic character

[1] R. Eichler *et al.*, RCA 98, 133 (2010),

[2] A. Yakushev *et al.*, Inorg. Chem. 53, 1624 (2014)

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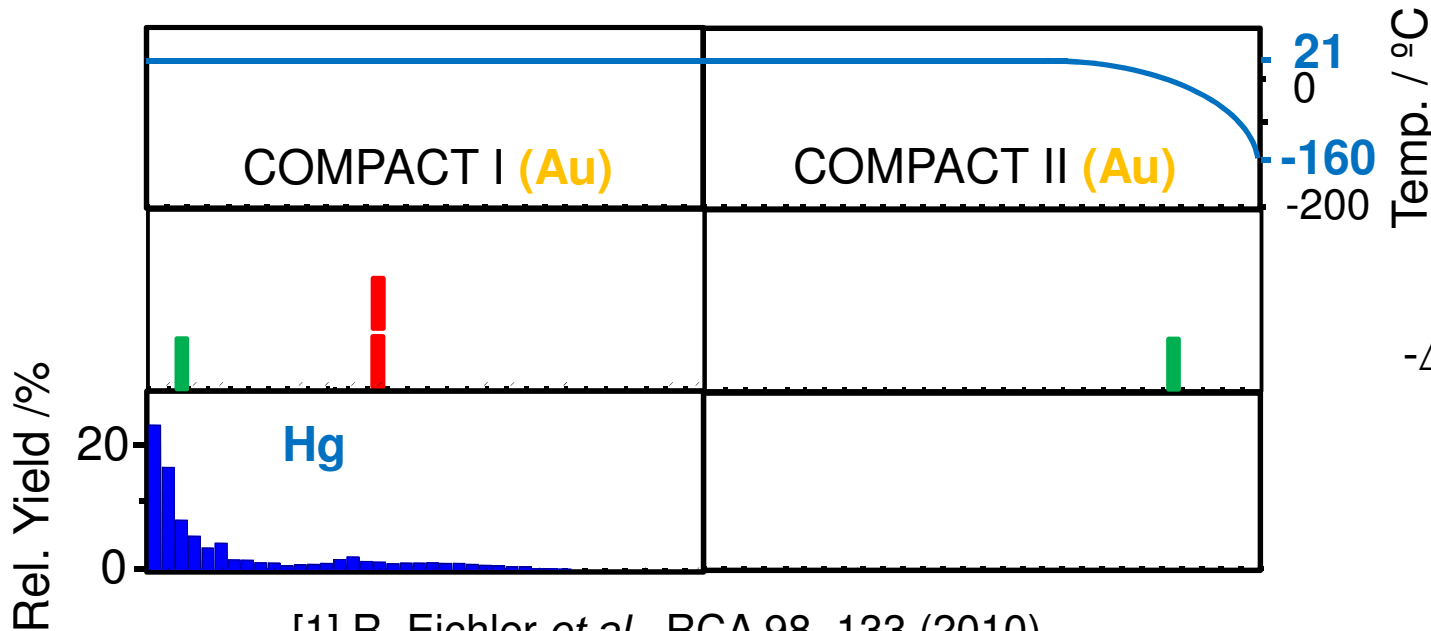


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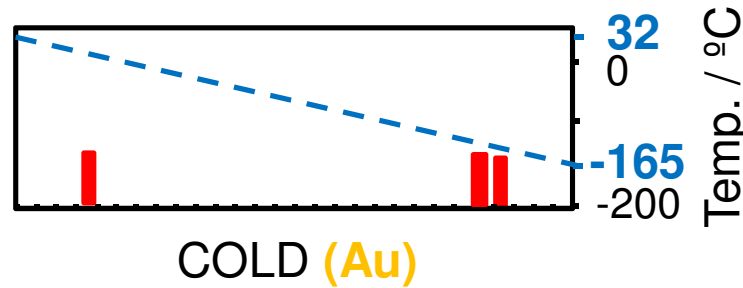
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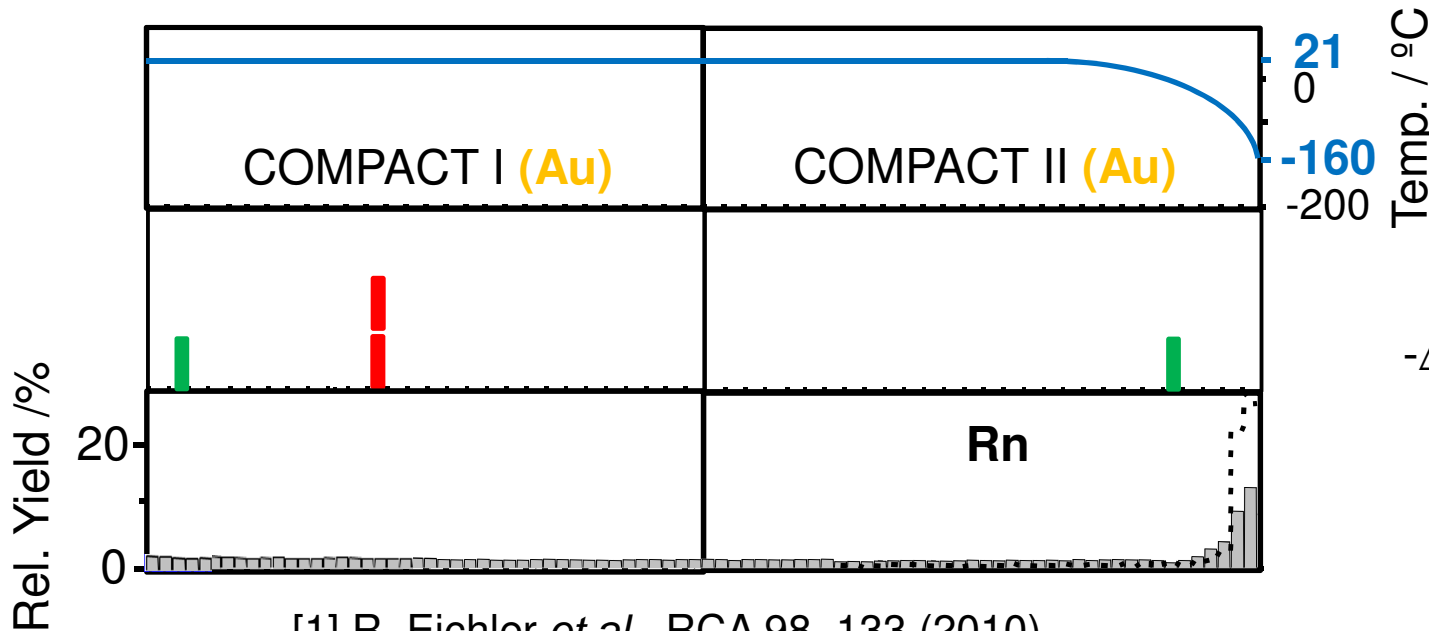


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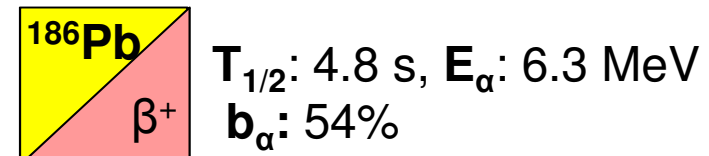
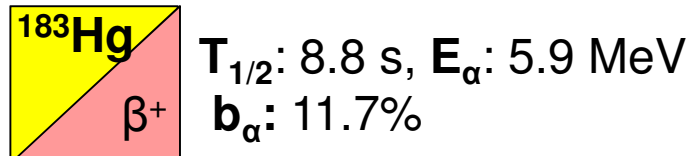
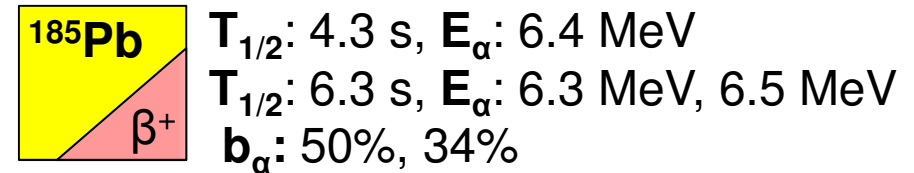
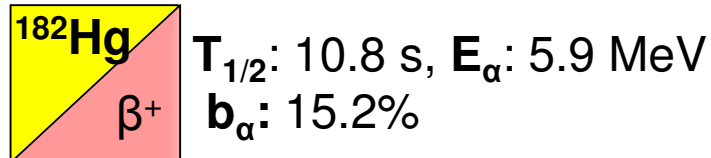
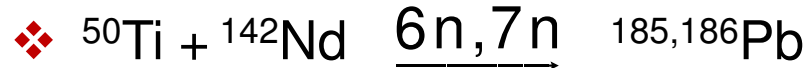
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[2] A. Yakushev *et al.*, Inorg. Chem. 53, 1624 (2014)

# Results of preparatory experiments with Pb and Hg

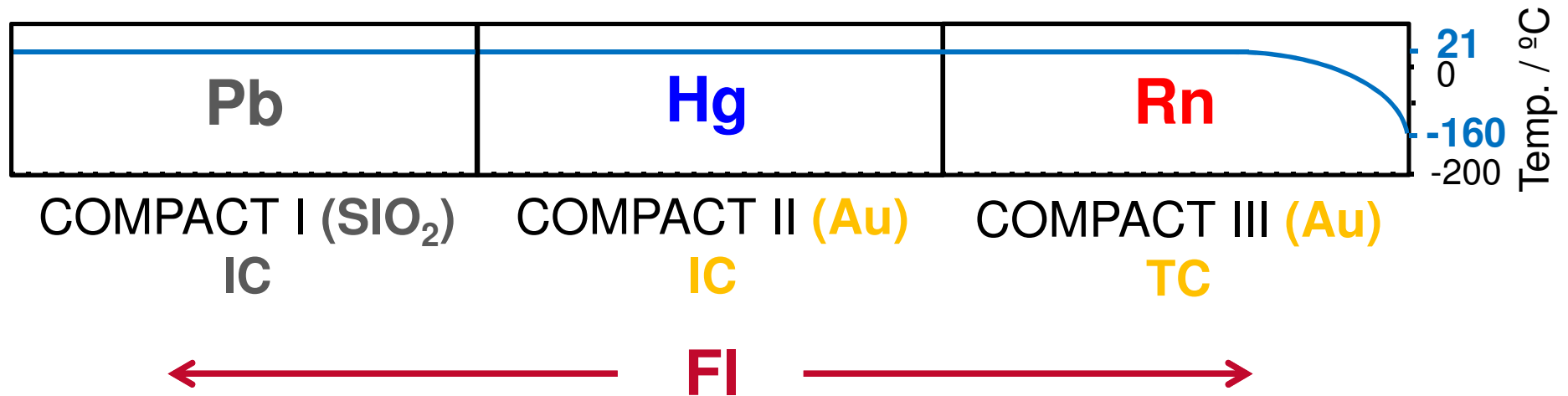
## Reaction



# FI experiment at TASCA 2014

## Scope:

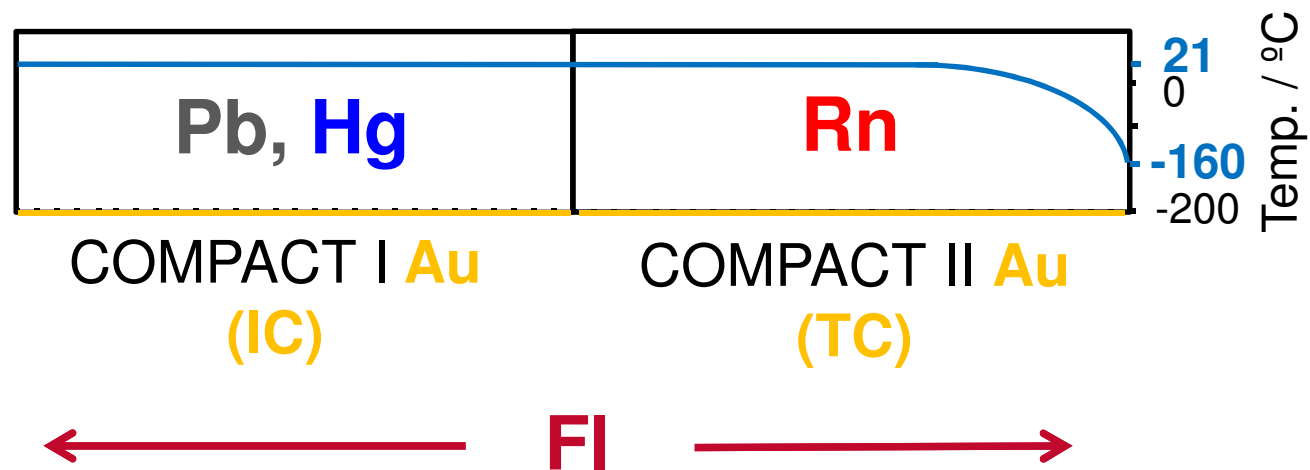
- ❖ Experimental investigations on the chemical properties of  $^{288,289}\text{Fl}$  in comparison to Pb, Hg, Rn and Cn



# The Flerovium Experiment 2015

## Scope:

- ❖ Optimization of the transport time with Hg
- ❖ Experimental investigations on the chemical properties of  $^{288,289}\text{Fl}$  in comparison to Pb, Hg, Rn and Cn





# Flerovium experiments 2014 + 2015

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Data are currently being analyzed