

# Non-target screening of new compounds in Arctic apex predators

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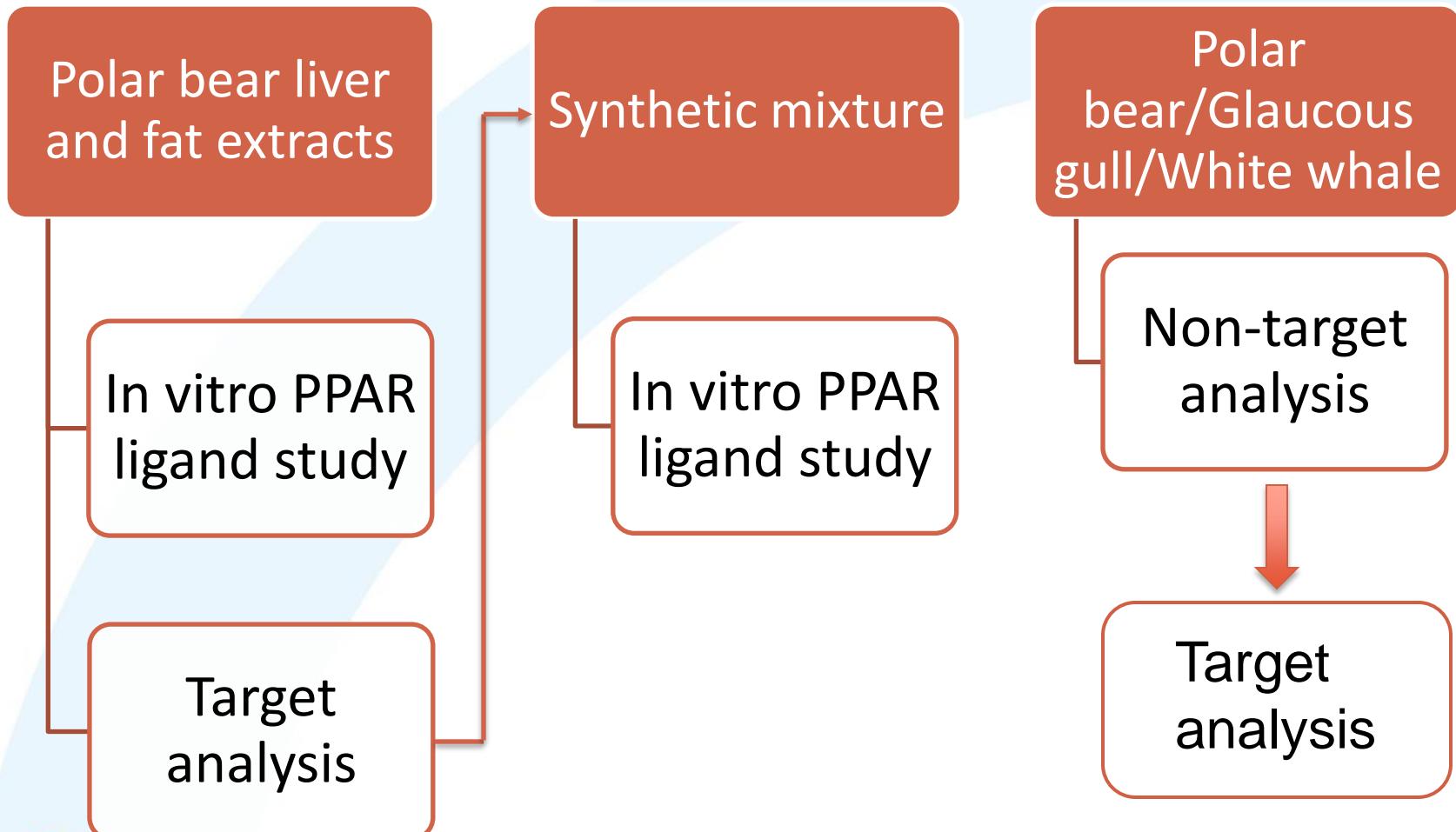
Norwegian Polar institute



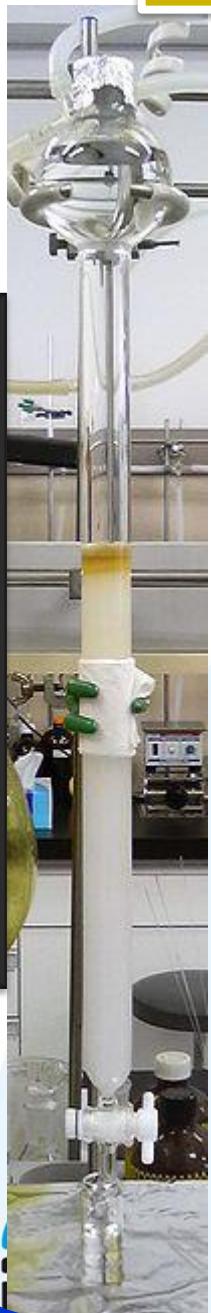
# Aim of the study

- The effects of Contaminants on polar bear energetics – using *in vitro* exposure studies on the peroxisome proliferator-activated receptors (PPARs)  
**Binding  $\longleftrightarrow$  Activation**
- Extracts as well as synthetic mixtures was compared
- Screening of potential new pollutants

# Project outline



## Extraction



68g Polarbear liver  
12g blubber  
Control/blank + QC

Homogenization  
Dry  $\text{Na}_2\text{SO}_4$

Extracting  
Lipids and POPs

Open-column Extraction

## Cleanup

Reverse Osmosis



Thin layflat polyethylene tubing  
 $<10 \text{ \AA}$  pores

Removes:  
88% lipids liver  
 $>98\%$  in blubber

GPC/SEC  
Size exclusion  
chromatography

Chromatography  
On Florisil

F1: Neutrals PCBs,  
PBDEs, pest etc  
F2:  $\text{MeSO}_2$ -PCBs,  
 $\text{MeSO}_2$ -DDE  
F3: OH-PCBs, PCP,  
OH-PBDEs

Conc factor 100 to 1

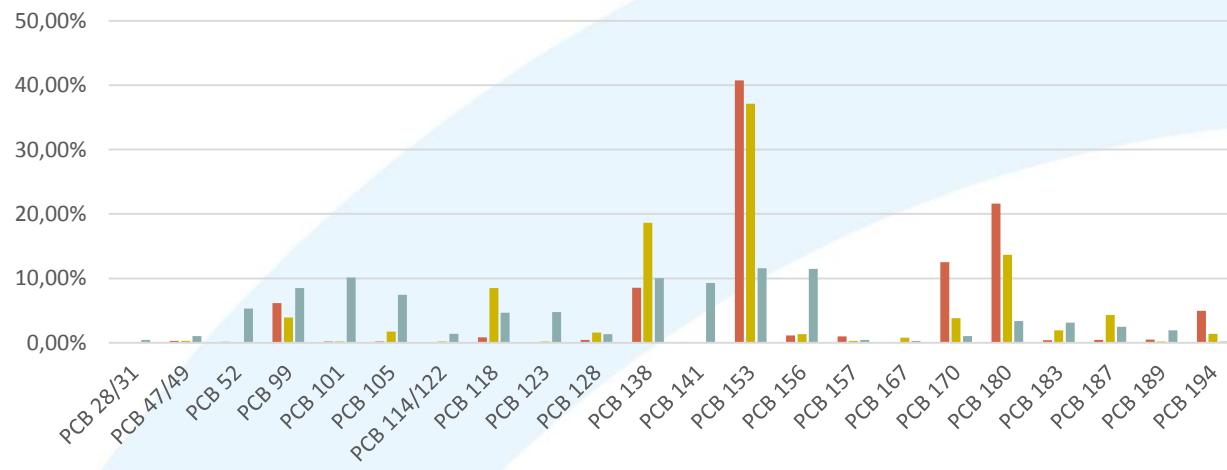
F1

F2

F3

# Target screening

PCB composition



Biotransformation :

■ Polar bear > ■ Glaucous gull > ■ White whale

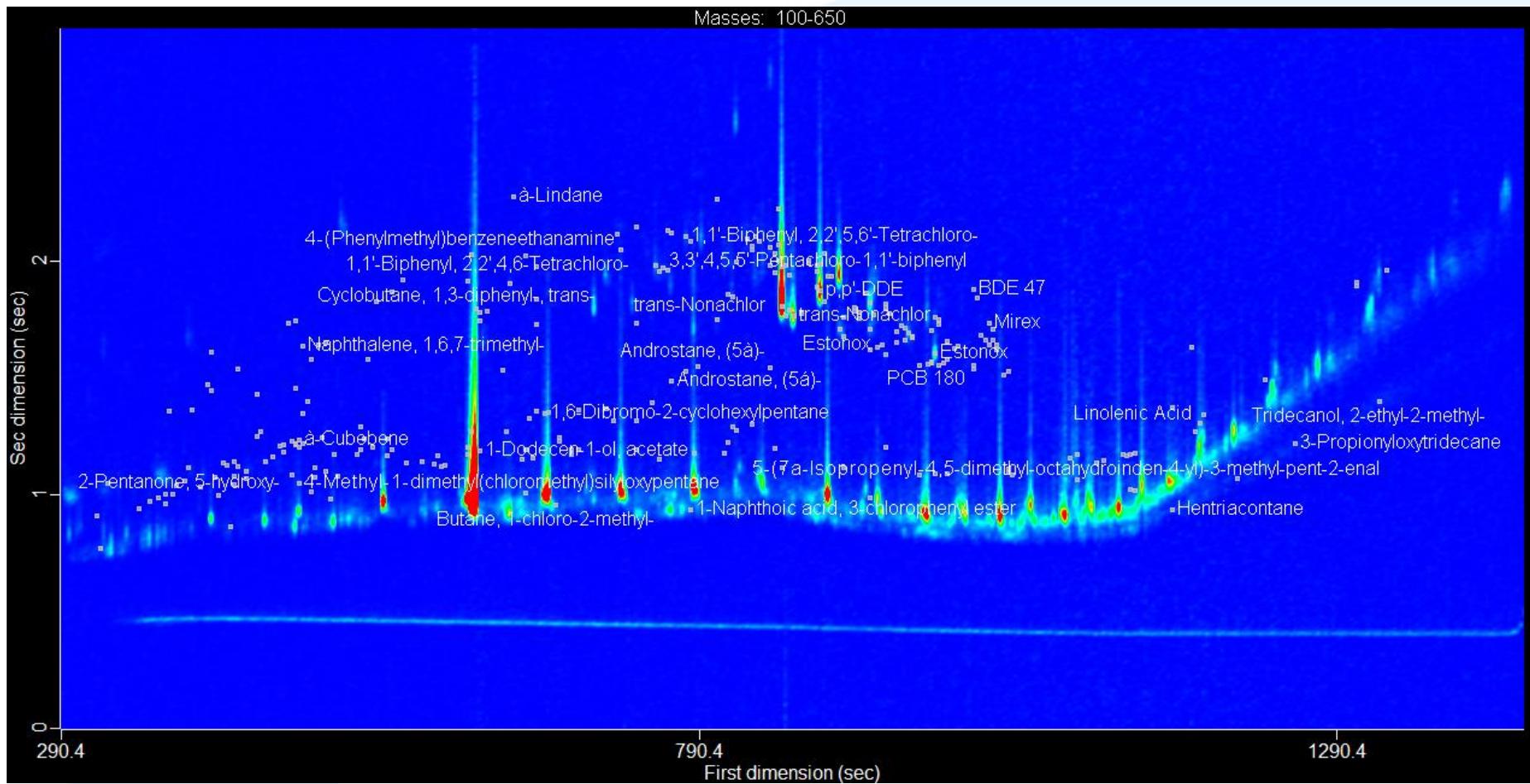


# Non-target analyses

- Fat samples from 3 polar bears, 3 glaucous gulls, 3 White whales
- Svalbard 2013-14
- Extraction: Homogenized in  $\text{Na}_2\text{SO}_4$  followed by extraction with acetone/cyclohexane
- Clean-up: semipermeable membrane, gel permeation chromatography and fractioning on Florisil to three according to polarity of the compounds
- Analyses: Leco Pegasus 4D, 2-D gas chromatografi coupled to time of flight mass spectrometry
- Data processing: Leco ChromaTOF software

Peak find → Peak deconvolution → matching to spectra database (NIST)  
→ scripting

# White whale



# Results of new compounds by non-target analyses

| Hit   | CAS         | Polar bear | Glaucous gull | White whale |
|---|-------------|------------|---------------|-------------|
| Octachlorostyrene   | 29082-74-4  | x          | x             |             |
| Heptachlor-endo-epoxide (isomer A)  | 28044-83-9  | x          | x             |             |
| 2,3-Dichlorobenz[b]thiophene  | 5323-97-7   | x          | x             |             |
| 1-Piperidinecarbodithioic acid, 2,3,5,6-tetrachloro-4-pyridinyl ester   |             | x          |               |             |
| (2,3-Diphenylcyclopropyl)methyl phenyl sulfoxide, trans-  | 131758-71-9 | x          |               | x           |
| 9-Allylanthracene   | 23707-65-5  |            |               | x           |
| L-Proline, n-pentafluoropropionyl-, heptyl ester  |             |            |               | x           |
| 2-Chloro-1,1-bis(4-chlorophenyl)-ethene (DDMU)  | 1022-22-6   |            | x             | x           |
| 7-Methylbenz(a)anthracene   | 2541-69-7   |            |               | x           |
| Tris(3-chlorophenyl)phosphine   | 29949-85-7  |            |               | x           |
| 6-Methoxy-2,2',4,4'-tetra (BC-3)  | 102739-99-1 |            |               | x           |
| Naphthalene, 1,2,3,4-tetrahydro-1-phenyl-   |             |            |               | x           |
| Bipyrrol Q1   | 428442-17-5 |            |               | x           |
| 1,3-Cyclopentadiene, 1,2,3,4-tetrachloro-   | 695-77-2    |            | x             |             |
| 1-Silacyclopentadiene, 3-(diethylboryl)-4-ethyl-1,1-dimethyl-2-(3-methylbutyl)-5-trimethylstannyl-                                | 153331-61-4 |            | x             |             |
| Endosulfan  | 115-29-7    |            | x             |             |
| 2,3,4,5,6-Pentachlorophenyl acetat  | 1441-02-7   |            | x             |             |
| 2,7:3,6-Dimethanonaphth[2,3-b]oxiren-8-ol, 3,4,5,6,9,9-hexachloro-1a,2,2a,3,6,6a,7,7a-octahydro-, acetate, stereoisomer           | 55649-85-9  |            | x             |             |
| 1,4:5,8-Dimethanonaphthalene-2,3-diol, 5,6,7,8,9,9-hexachloro-1,2,3,4,4a,5,8,8a-octahydro-, diacetate, (1a,2a,3a,4a,4a,5a,8a,8a)- | 34408-22-5  |            | x             |             |
| Chlordecone   | 55570-85-9  |            | x             |             |

- Qualitative
- No information of concentrations
- Structures not confirmed by standards

# Target analysis

- Fat samples from 10 polar bears, 8 glaucous gulls, 9 white whales
- Extraction and cleanup similar as non-target analysis (no semipermeable membrane)– full set of internal standards
- Analysis on GC-MS

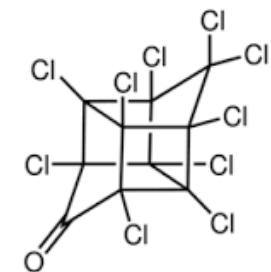
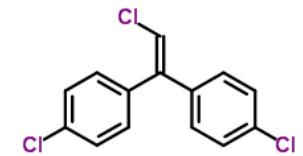
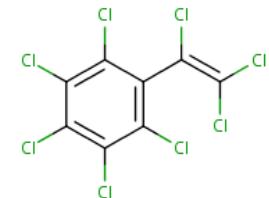
# Target analyses of compounds found in non-target screening:

| Concentrations (ng/g ww)  | Polar bear       | Glaucous gull | White whale |
|---|------------------|---------------|-------------|
| Octachlorostyrene   | 8.7              | 21.6          | 3.8         |
| Heptachlor-endo-epoxide   | < 0.9            | < 0.9         | < 0.9       |
| 2,3-Dichlorobenzo[b]thiophene   | < 0.1            | < 0.1         | < 0.1       |
| 1-Piperidinecarbodithioic acid, 2,3,5,6-tetrachloro-4-pyridinyl ester   |                  |               |             |
| (2,3-Diphenylcyclopropyl)methyl phenyl sulfoxide, trans-  |                  |               |             |
| 9-Allylanthracene   |                  |               |             |
| I-Proline, n-pentafluoropropionyl-, heptyl ester  |                  |               |             |
| 2-Chloro-1,1-bis(4-chlorophenyl)-ethene (DDMU)  | 2.5              | 54.3          | 3.5         |
| 7-Methylbenz(a)anthracene   | < 0.1            | < 0.1         | < 0.1       |
| Tris(3-chlorophenyl)phosphine   | < 0.3            | < 0.3         | < 0.3       |
| 6-Methoxy-2,2',4,4'-tetra (BC-3)  |                  |               |             |
| Naphthalene, 1,2,3,4-tetrahydro-1-phenyl-   | < 2.5            | < 2.5         | < 2.5       |
| Bipyrrol Q1   |                  |               |             |
| 1,3-Cyclopentadiene, 1,2,3,4-tetrachloro-   | < 0.3            | < 0.3         | < 0.3       |
| 1-Silacyclopentadiene, 3-(diethylboryl)-4-ethyl-1,1-dimethyl-2-(3-methylbutyl)-5-trimethylstannyl-                                |                  |               |             |
| Endosulfan  | Will be analyzed |               |             |
| 2,3,4,5,6-Pentachlorophenyl acetat  |                  |               |             |
| 2,7:3,6-Dimethanonaphth[2,3-b]oxiren-8-ol, 3,4,5,6,9,9-hexachloro-1a,2,2a,3,6,6a,7,7a-octahydro-, acetate, stereoisomer           |                  |               |             |
| 1,4:5,8-Dimethanonaphthalene-2,3-diol, 5,6,7,8,9,9-hexachloro-1,2,3,4,4a,5,8,8a-octahydro-, diacetate, (1a,2a,3a,4a,4a,5a,8a,8a)- |                  |               |             |
| Chlordecone   | < 0.1-2.5        | 24.7          | 1.7         |

- New extractions
- Quantification using reference standards
- Retention time confirmed using authentic standards
- Quantitative results

# Target analyses of compounds found in non-target screening:

- Octachlorostyrene:
  - Industrial by-product, not regulated
  - Not new, reported in polar bears already in 2000
  - Highly toxic
- DDMU:
  - Metabolite of DDE
  - Toxic
- Chlordcone/Kepone
  - Pesticide, included in Stockholm Convention in 2009
  - Not measured in the Arctic (Muir & deWit, 2010)
  - Toxic



# The way forward

- Suspect screening, good alternative or complement to non-target screening
- Accurate high resolution MS for additional confirmation - standards are difficult and expensive to get
- Confirmation of compounds – Retention time, accurate mass, mass fragmentation spectra, isotope ratios
- Negative and positive chemical ionization  
→Selectivity and molecular fragments

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