

PERSISTENT ORGANIC POLLUTANTS IN THE LIVERS OF MOOSE HARVESTED IN THE SOUTHERN NORTHWEST TERRITORIES, CANADA

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ANALYTICAL METHODS PCBs, OCOs, and BFRs

Samples were thawed and thoroughly homogenized in a small stainless steel blender. Subsamples (3.5 to 5 g ww) were Soxhlet-extracted with dichloromethane; all were spiked with 1,3,5-tribromobenzene (TrBB) prior to extraction. Recovery of brominated compounds was monitored using BDE-71, d16-gHBCD, and C13-BDE-209. ¹³C₁₂-PCB-133 was added as a recovery standard for GPC performance for samples analyzed by ALS Global. A laboratory blank consisting of all reagents and 2 NIST reference materials (cod liver SRM 1588b and fish muscle SRM 1946) were also analyzed with the 14 samples. The extracts were then rotary evaporated under vacuum, exchanged into DCM:hexane (1:1) and applied to a gel permeation chromatography (GPC) column (60 g Bio-beads SX3) to remove lipids and other biogenic materials. The GPC eluate was reduced to 1 mL under vacuum. Percent lipid was determined gravimetrically on a sub-sample of the extract or by evaporating the first GPC fraction.

Extracts were cleaned up on a silica gel column. NLET utilized activated silica gel (8 g, 1.1 cm i.d. column), and eluted with hexane followed by *n*-hexane/DCM (1:1) to separate PCBs from most of the PBDEs

and OCPs. BDE-209 was quantitatively eluted in the silica Fraction 1.

The ALS Global split the GPC elution into separate OCP/OCO fractions that were chromatographed on a 2% deactivated silica gel column and then reduced to 0.05 mL for analysis. The PCB fraction was cleaned up on an acid-silica gel column (45% w/w H₂SO₄ on Silica Gel topped with neutral Silica Gel) then reduced to 0.04 mL for analysis.

GC-ECD analysis was conducted on 7 Dehcho samples using a GC-ECD (Agilent 6890 gas chromatograph with a ⁶³Ni-electron capture detector [ECD]) using a 30 m x 0.25 mm (i.d.) DB-5 column (internal film thickness 0.25 mm; J&W Scientific, Folsom, California, USA) with H₂ carrier gas (constant flow rate 0.91 mL min⁻¹). Ultra-pure N₂ was used as the makeup gas for the ECD (detector temperature: 325 C). The GC-ECD quantification of OCs in each sample was performed using a 4-point external standard calibration curve. Calibration standards were quantified after every 10 samples.

Toxaphene-related compounds, including 22 polychlorinated bornane congeners as well as α - and β -endosulfan and endosulfan sulfate, were quantified by GC-electron capture-negative ion mode (ECNI) mass spectrometry using an Agilent 6890 GC-5975

MS system as described by Hoekstra et al. (2002). Toxaphene and homologues were quantified using a “Hercules” technical standard as described by Glassmeyer et al. (1999). Individual polychlorinated bornane congeners were quantified using a series of external calibration standards (Dr. Ehrenstofer, Augsburg, Germany). Alpha- and beta-endosulfan, and endosulfan sulfate were quantified by external standards using the characteristic fragment ions m/z 406 and m/z 273.

Analyses of all PBDEs and other brominated flame retardants (BFRs) was carried out by GC-ECNIMS on an Agilent 6890-5975 MS using an HP5-MS capillary column (30 m x 0.25 mm x 0.25 um film thickness). Helium was the carrier gas, and separation was performed at a constant flow of 1.2 mL/min (Muir et al. 2006). The mass spectrometer was operated in the NCI mode, methane was the buffer gas, and temperature was 106, 150, and 300 °C for the quadrupole, the ion source, and the interface, respectively. The analytes were monitored at m/z 79/81 using an external standard calibration, except for C13-BDE-209 which was monitored at m/z 493/495 and native BDE-209 at m/z 487/485. Any β - and γ -HBCDD residues in the samples were most likely thermally isomerized to α -HBCDD in the GC injection port, thus, results represent total HBCDD (Muir et al. 2006).

PFASs

An internal standard mixture of ^{13}C -PFASs was added to every sample and extracted by shaking twice with acetonitrile. The extract was evaporated under nitrogen to dryness and reconstituted with 1 mL of methanol. The extract was cleaned with a graphite

carbon solid phase cartridge (Supelco). Cleaned up extracts were analyzed for PFCAs as well as PFSAs. The analyses were performed by liquid chromatography with negative electrospray tandem mass spectrometry (LC-MS/MS). Analytes were detected using an API 4000 Q Trap (Applied Biosystems, Carlsbad, California, USA) after chromatographic separation with an Agilent 1100 LC. Chromatography was performed using an ACE C18 column (50 mm x 2.1 mm, 3 μm particle size; Aberdeen, United Kingdom), preceded by a C18 guard column (4.0 x 2.0 mm, Phenomenex) and the column oven was set to 30 °C. Samples were quantified with a 6 point calibration curve and isotopic dilution method.

QUALITY ASSURANCE

Recoveries of internal standards ranged from 79% for δ -HCH to 120% for PCB-204 (Table S4) in Decho samples analysed by GC-ECD, and from 59% for endrin ketone to 138% for 1245-TTBB in samples analysed by GC-HRMS/LRMS. A recovery spike demonstrated good recoveries of 32 OCP/OCOs (55–101%) and 18 PBDE/BFRs (71–133%) (Table S5). Slight losses of more volatile compounds (e.g., hexachlorobutadiene) and recovery enhancement due to contribution from laboratory reagent blanks (BDE 47) explain the recovery variation. No corrections for recovery were made based on this information. Analysis of the reference materials (NIST SRMs 1588b and 1946) showed good agreement with all analytes quantified to within $\pm 25\%$ of certified values of OCP/OCOs (17 compounds) and PCBs (29 congeners).

Table S1. List of moose samples, collection year, age, sex, biological characteristics, harvest date and location.

Region	ID	Age	Sex	Condition E/G/F ¹	Hunter est. age	Date Harvested	Latitude °N	Longitude °W
South Slave	SSR-MO-11-4	<1	M	F	Calf	20-Feb-10	60.21	112.13
	SSR-MO-11-7	1	M	E	Adult	10-Jan-10	60.64	112.27
	SSR-MO-11-8	4	M	n/a	Adult	4-Oct-10	60.76	112.19
	SSR-MO-11-9	4	M	G	Adult	13-Dec-10	60.79	116.22
	SSR-MO-11-10	<1	M	G	Yearling	30-Sep-10	60.85	114.49
	SSR-MO-11-12	<1	F	G	calf	8-Feb-10	60.78	112.78
	SSR-MO-11-13	9	F	n/a	Adult	8-Feb-10	60.78	112.78
Dehcho	GEN-14	3	M	G	Adult	21-Feb-06	61.55	121.18
	GEN-15	6	F	G	Adult	02-Mar-06	61.64	121.06
	GEN-16	12	F	G	Adult	06-Mar-06	61.60	121.14
	GEN-17	<1	F	G	Calf	06-Mar-06	61.60	121.14
	GEN-18	<1	M	G	Calf	24-Mar-06	61.50	120.62
	GEN-19	2	M	E	Adult	18-Mar-06	61.48	120.62
	JMR-4	5	F	E	Adult	24-Mar-06	61.49	120.64

¹ Hunter condition evaluation. E = excellent, G= good, F = fair. n/a = not available.

Table S2. List of individual organohalogen analytes along with their MDLs (ng/g ww) using either GC with high or low resolution MS, GC-NCIMS, or GC-ECD.

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	MDL
				GC-MS	GC-ECD
OCO	Hexachlorobutadiene		GC-HRMS, GC-ECD	<0.005	0.032
OCO	1,2,4,5-Tetrachlorobenzene	1,2,4,5-tetrachlorobenzene	GC-HRMS, GC-ECD	0.099	0.007
OCO	1,2,3,4-Tetrachlorobenzene	1,2,3,4-tetrachlorobenzene	GC-HRMS, GC-ECD	0.108	0.015
OCP	Pentachlorobenzene	pentachlorobenzene	GC-HRMS, GC-ECD	0.077	<0.002
OCP	345-trichlorovertatrole		GC-HRMS, GC-ECD	na	0.012
OCP	Pentachloroanisole		GC-HRMS, GC-ECD	<0.008	0.024
OCP	Hexachlorobenzene		GC-HRMS, GC-ECD	<0.026	<0.002
OCP	3,4,5,6-Tetrachlorovertatrole	GC-HRMS, GC-ECD	<0.016	<0.011	
OCP	α -HCH	α -hexachlorocyclohexane	GC-HRMS, GC-ECD	<0.059	0.015
OCP	β -HCH	β -hexachlorocyclohexane	GC-HRMS, GC-ECD	<0.1	0.013
OCP	γ -HCH	lindane	GC-HRMS, GC-ECD	<0.075	<0.002
OCP	Heptachlor		GC-HRMS, GC-ECD	<0.018	0.08
OCP	Pentachloronitrobenzene		GC-HRMS, GC-ECD	<0.075	na
OCP	Aldrin		GC-HRMS, GC-ECD	<0.011	0.026
OCP	Dacthal		GC-HRMS, GC-ECD	<0.056	na
OCP	Octachlorostyrene		GC-HRMS, GC-ECD	<0.025	0.016
OCP	HeptachlorEpoxide		GC-HRMS, GC-ECD	<0.069	0.047
OCP	Oxychlordane		GC-HRMS, GC-ECD	<0.019	<0.002
OCP	trans-chlordane		GC-HRMS, GC-ECD	<0.037	0.005
OCP	cis-chlordane		GC-HRMS, GC-ECD	<0.032	0.012
OCP	trans-nonachlor		GC-HRMS, GC-ECD	<0.035	0.062
OCP	Dieldrin		GC-HRMS, GC-ECD	<0.018	0.088
OCP	cis-nonachlor		GC-HRMS, GC-ECD	<0.039	0.024
OCP	Endrin		GC-HRMS, GC-ECD	<0.387	0.023

Table S2 continued . . .

Table S2 continued

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	MDL
				GC-MS	GC-ECD
OCP	24'-DDE	GC-HRMS, GC-ECD	<0.01	0.001	
OCP	44'-DDE	GC-HRMS, GC-ECD	<0.01	0.021	
OCP	24'-DDD	GC-HRMS, GC-ECD	<0.01	0.014	
OCP	44'-DDD	GC-HRMS, GC-ECD	<0.01	0.02	
OCP	24'-DDT	GC-HRMS, GC-ECD	<0.01	0.02	
OCP	44'-DDT	GC-HRMS, GC-ECD	<0.01	0.02	
OCP	Methoxychlor	GC-HRMS, GC-ECD	<0.01	0.019	
OCP	Mirex	GC-HRMS, GC-ECD	0.068	0.01	
PCBs	PCB-1	GC-LRMS, GC-ECD	<0.002	na	
PCBs	PCB-3	GC-LRMS, GC-ECD	<0.002	na	
PCBs	PCB4/10	GC-LRMS, GC-ECD	<0.002	0.166	
PCBs	PCB7/9	GC-LRMS, GC-ECD	<0.002	0.158	
PCBs	PCB6	GC-LRMS, GC-ECD	<0.002	0.042	
PCBs	PCB8/5	GC-LRMS, GC-ECD	<0.002	0.078	
PCBs	PCB12/13	GC-LRMS, GC-ECD	<0.002	1.11	
PCBs	PCB15	GC-LRMS, GC-ECD	<0.002	0.036	
PCBs	PCB19	GC-LRMS, GC-ECD	<0.002	<0.002	
PCBs	PCB18	GC-LRMS, GC-ECD	<0.002	0.211	
PCBs	PCB17	GC-LRMS, GC-ECD	<0.002	na	
PCBs	PCB27/24	GC-LRMS, GC-ECD	<0.002	0.036	
PCBs	PCB16/32	GC-LRMS, GC-ECD	<0.002	0.119	
PCBs	PCB26	GC-LRMS, GC-ECD	<0.002	0.112	
PCBs	PCB25	GC-LRMS, GC-ECD	<0.002	0.019	

Table S2 continued . . .

Table S2 continued

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	MDL
				GC-MS	GC-ECD
PCBs	PCB31	trichlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.445
PCBs	PCB28	trichlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.05
PCBs	PCB20/33/21	trichlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.155
PCBs	PCB22	trichlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.097
PCBs	PCB37	trichlorobiphenyl	GC-LRMS, GC-ECD	<0.002	
PCBs	PCB53	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.117
PCBs	PCB45	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.136
PCBs	PCB46	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.058
PCBs	PCB73/52	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.913
PCBs	PCB43/49	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.446
PCBs	PCB48/47/75	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	5.152
PCBs	PCB44	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.554
PCBs	PCB59/42	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.21
PCBs	PCB71/41/68/64	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.417
PCBs	PCB100	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.052
PCBs	PCB63	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.024
PCBs	PCB74/61	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.241
PCBs	PCB70/76	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.663
PCBs	PCB80/66	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.375
PCBs	PCB56/60	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.229
PCBs	PCB81	tetrachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.413
PCBs	PCB95/93	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.641
PCBs	PCB91	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.146

Table S2 continued . . .

Table S2 continued

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	MDL
				GC-MS	GC-ECD
PCBs	PCB92	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.156
PCBs	PCB84/90/01/89	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.26
PCBs	PCB89-101	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.692
PCBs	PCB99	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.319
PCBs	PCB119	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.022
PCBs	PCB83/108	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.089
PCBs	PCB97	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.172
PCBs	PCB86/111/125/117/87/116/115	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	na
PCBs	PCB120/85	pentachlorobiphenyl	GC-LRMS, GC-ECD	0.055	0.092
PCBs	PCB110	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.461
PCBs	PCB136	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.114
PCBs	PCB82	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.049
PCBs	PCB107/109	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.059
PCBs	PCB123	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.406
PCBs	PCB118/106	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.298
PCBs	PCB114	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.024
PCBs	PCB105/127	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.141
PCBs	PCB126	pentachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	na
PCBs	PCB151	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.173
PCBs	PCB135/144	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.651
PCBs	PCB139/149	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	na
PCBs	PCB131/165/142	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.021
PCBs	PCB146	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.056

Table S2 continued . . .

Table S2 continued

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	MDL
				GC-MS	GC-ECD
PCBs	PCB153	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.293
PCBs	PCB132/168	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.135
PCBs	PCB141	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.068
PCBs	PCB137	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	3.971
PCBs	PCB163/164/138	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.283
PCBs	PCB158/160	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.044
PCBs	PCB129	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.018
PCBs	PCB159	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	na
PCBs	PCB128	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.028
PCBs	PCB167	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.01
PCBs	PCB156	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.013
PCBs	PCB157	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	<0.001
PCBs	PCB169	hexachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	na
PCBs	PCB182/187	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.109
PCBs	PCB183	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.046
PCBs	PCB174/181	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.072
PCBs	PCB177	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.143
PCBs	PCB171	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.042
PCBs	PCB172/192	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.018
PCBs	PCB197	octachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.007
PCBs	PCB180	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.087
PCBs	PCB193	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.132
PCBs	PCB191	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	<0.001

Table S2 continued . . .

Table S2 continued

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	MDL
				GC-MS	GC-ECD
PCBs	PCB170/190	heptachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.03
PCBs	PCB-202	octachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	na
PCBs	PCB199	octachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.008
PCBs	PCB196/203	octachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.022
PCBs	PCB195	octachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.015
PCBs	PCB194	octachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.005
PCBs	PCB205	octachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	<0.001
PCBs	PCB208	nonachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.002
PCBs	PCB207	nonachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	<0.001
PCBs	PCB206	nonachlorobiphenyl	GC-LRMS, GC-ECD	<0.002	0.013
PCBs	PCB209	decachlorobiphenyl	GC-LRMS, GC-ECD	0.08	0.004
Sum op- and pp'-DDT related					
Σ DDT					
Σ CHL					
Sum oxychlordane, C & t-chlordane, c & t-nonachlor, heptachlor, heptachlor epoxide					
Sum α , β , γ -HCH					
Σ HCH					
Sum tetra, PeCBz, HCB					
Σ CBz					
Sum endosulfan					
Σ endosulfan					
Sum PCBs					
Σ PCB					
Sum mono-di CBs					
Σ mono-di					
Sum tri CBs					
Σ tri					
Sum tetra cBs					
Σ tetra					
Sum penta CBs					
Σ penta					
Sum hexa CBs					
Σ hexa					

Table S2 continued . . .

Table S2 continued

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	MDL
				GC-MS	GC-ECD
Sum hepta CBs	Σ hepta				
Sum octa CBs	Σ octa				
Sum nona-deca CBs	Σ nona-deca			<0.002	
Endosulfan	α -Endosulfan			<0.002	
Endosulfan	β -Endosulfan			<0.002	
Endosulfan	Endosulfan sulfate			<0.002	
Perfluorosulfonates	PFBS	Perfluorobutane sulfonate	LC-MS/MS	<0.001	
Perfluorosulfonates	PFHxS	Perfluorohexane sulfonate	LC-MS/MS	<0.001	
Perfluorosulfonates	PFHpS	Perfluoroheptane sulfonate	LC-MS/MS	<0.001	
Perfluorosulfonates	PFOS	Perfluorooctane sulfonate	LC-MS/MS	0.299	
Perfluorosulfonates	PFDS	Perfluorodecane sulfonate	LC-MS/MS	<0.001	
Perfluorosulfonates	PFOSA	Perfluoroctanesulfonamide	LC-MS/MS	<0.001	
Perfluorosulfonates	Σ PFSA _s	Sum perfluorosulfonates	LC-MS/MS		
Perfluorocarboxylates	PFHxA	Perfluorohexanoate	LC-MS/MS	0.179	
Perfluorocarboxylates	PFHpA	Perfluoroheptanoate	LC-MS/MS	0.036	
Perfluorocarboxylates	PFOA	Perfluorooctanoate	LC-MS/MS	0.213	
Perfluorocarboxylates	PFNA	Perfluorononanoate	LC-MS/MS	0.046	
Perfluorocarboxylates	PFDA	Perfluorodecanoate	LC-MS/MS	0.256	
Perfluorocarboxylates	PFUnA	Perfluoroundecanoate	LC-MS/MS	0.085	
Perfluorocarboxylates	PFDoA	Perfluoroundodecanoate	LC-MS/MS	0.159	
Perfluorocarboxylates	PFTA	Perfluorotridecanoate	LC-MS/MS	0.089	
Perfluorocarboxylates	PFTnA	Perfluorotetradecanoate	LC-MS/MS	0.131	
Perfluorocarboxylates	Σ PFCA _s	Sum perfluorocarboxylates	LC-MS/MS		

Table S2 continued . . .

Table S2 continued

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	GC-MS	MDL
						GC-ECD
Toxaphene	Toxaphene	Technical standard	GC-ECNIMS	<0.05		
Toxaphene homologs	Hexa	Technical standard	GC-ECNIMS	<0.05		
Toxaphene homologs	Hepta	Technical standard	GC-ECNIMS	<0.05		
Toxaphene homologs	Octa	Technical standard	GC-ECNIMS	<0.05		
Toxaphene homologs	Nona	Technical standard	GC-ECNIMS	<0.05		
Toxaphene homologs	Deca	Technical standard	GC-ECNIMS	<0.05		
Chlorobornanes ²	Parlar 11-12	Technical standard	GC-ECNIMS	<0.05		
Chlorobornanes	Parlar 15		GC-ECNIMS	<0.02		
Chlorobornanes	Hex-sed		GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 21		GC-ECNIMS	<0.02		
Chlorobornanes	Hep-sed		GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 25		GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 32	B7-499	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 26	B7-515	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 31	B8-1413	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 38	B8-789	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 39	B8-531	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 40-41	B8-1414/1945	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 42	B8-806	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 44	B8-2229	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 50	B9-1679	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 51	B8-786	GC-ECNIMS	<0.02		
Chlorobornanes	Parlar 56	B9-1046	GC-ECNIMS	<0.02		

Table S2 continued . . .

Table S2 continued

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	MDL
				GC-MS	GC-ECD
Chlorobornanes	Parlar 58	B9-715	GC-ECNIMS	<0.02	
Chlorobornanes	Parlar 59	B9-1049	GC-ECNIMS	<0.02	
Chlorobornanes	Parlar 62	B9-1025	GC-ECNIMS	<0.02	
Chlorobornanes	Parlar 63	B9-2206	GC-ECNIMS	<0.02	
Chlorobornanes	Parlar 69	B10-1110	GC-ECNIMS	<0.02	
PBDEs	BDE 17	dibromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 28/33	tribromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 49	tetrabromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 47	tertetrabromodiphenyl ether	GC-ECNIMS	<0.035	
PBDEs	BDE 66	terabromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 100	pentabromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 99	pentabromodiphenyl ether	GC-ECNIMS	<0.022	
PBDEs	BDE 85	pentabromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 154	hexabromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 153	hexabromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 138	hexabromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 183	heptabromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 190	heptabromodiphenyl ether	GC-ECNIMS	<0.002	
PBDEs	BDE 209	decabromodiphenyl ether	GC-ECNIMS	<0.558	
Sum PBDEs					
Other BFRs	TBP-AE	Allyl 2,4,6-tribromophenyl ether	GC-ECNIMS	<0.002	
Other BFRs	pTBX	tetrabromoxylene	GC-ECNIMS	<0.002	
Other BFRs	TBP-DBPE	2-Bromoallyl 2,4,6-tribromophenyl ether	GC-ECNIMS	<0.002	

Table S2 continued . . .

Table S2 continued

Class ¹	Common name	Chemical or other name	Instrumental analysis	MDL ³	GC-MS	MDL
						GC-ECD
Other BFRs	PBB _e	pentabromobenzene	GC-ECNIMS	<0.002		
	TBC _T	terabromo-o-chlorotoluene	GC-ECNIMS	<0.002		
Other BFRs	PBTo	pentabromotoluene	GC-ECNIMS	<0.002		
Other BFRs	PBEB	pentabromoethylbenzene	GC-ECNIMS	<0.002		
Other BFRs	DPTE/TBP-DBP _E	2,3-Dibromopropyl 2,4,6-tribromophenyl ether	GC-ECNIMS	<0.002		
Other BFRs	HBB	hexabromobenzene	GC-ECNIMS	<0.002		
Other BFRs	BB-101	penetabromobiphenyl	GC-ECNIMS	<0.002		
Other BFRs	PBBA	pentabromobenzyl acrylate	GC-ECNIMS	<0.002		
Other BFRs	EHTeBB	2-ethyl-1-hexyl 2,3,4,5-tetrabromobenzoate	GC-ECNIMS	<0.002		
Other BFRs	HBCDD	hexabromocyclododecane	GC-ECNIMS	<0.002		
Other BFRs	BTBPE	bis(tribromophenoxy) ethane	GC-ECNIMS	<0.002		
Other BFRs	BEHTBP	Bis(2-ethyl-1-hexyl) tetrabromophthalate	GC-ECNIMS	<0.002		
OCO	syn-DP	Syn-Dechlorane	GC-ECNIMS	<0.002		
OCO	anti-DP	Anti-Dechlorane	GC-ECNIMS	<0.002		
Other BFRs	OBIND	Octabromotrimethylphenylindane	GC-ECNIMS	<0.002		

¹OCO = other chlorinated organic, OCP = organochlorine pesticide related, PCB = polychlorinated biphenyl, PBDE = polybrominated diphenyl ether, BFR = brominated flame retardant.

²Nomenclature of chlorobornanes based on Andrews and Vetter 1995.

³MDL = method detection limit = 3*SD of blank values. “na” = not analysed. Where nondetect values were present in blanks a “<” values were used. These are instrument detection limits based on S/N of approximately 10:1.

Table S3. Comparison of GC-MS and GC-ECD analysis of 3 moose liver samples. See Table S2 for full list analytes represented by each group.

Target Analytes	moose	moose	moose	moose	moose	moose
	GC-MS	ECD	GC-MS	ECD	GC-MS	ECD
	GEN-14	Gen-14	GEN-15	Gen-15r	GEN-16	Gen-16
ng/g ww	ng/g ww	ng/g ww	ng/g ww	ng/g ww	ng/g ww	ng/g ww
ΣDDT	<0.002	<0.002	<0.002	<0.002	<0.002	0.01
ΣCHL	0.09	0.02	0.02	0.05	0.02	0.02
ΣHCH	0.11	0.13	0.07	0.11	0.07	0.12
HCB	0.19	0.17	0.24	0.12	0.20	0.15
ΣPCB	0.62	0.30	1.07	0.74	1.03	0.89
smono-di	0.22	0.06	0.44	0.11	0.21	0.13
Σ-tri	0.26	0.13	0.11	0.15	0.11	0.31
Σ-tetra	0.08	0.08	0.33	0.19	0.42	0.27
Σ-penta	0.05	<0.002	0.18	0.17	0.06	0.02
Σ-hexa	0.02	0.01	0.02	0.03	0.02	0.07
Σ-endosulfan	0.01	0.07	0.03	0.04	0.03	0.12

Table S4. Recoveries of internal standards during or prior to sample extraction.

Compound	% recovery		% recovery	
	GC-ECD analysis (n = 7)	SD	GC-MS analysis (n = 10)	SD
1,3-DBB	84.7	3.5	85.6	23.9
1,3,5-TBB	80.2	3.8	112	29.2
1,2,4,5-TTBB	87.7	4.3	138	35.3
δ-HCH	79.0	5.2	110	39.3
Endrin Ketone	91.8	5.1	59.4	37.2
PCB 30	108	3.3	94.0	5.1
PCB 204	120	4.4	98.6	3.0
D16-gHBCDD ¹			84.1	17.7
¹³ C-BDE-209 ¹			67.4	15.8
¹³ C12-PCB133			82.3	8.0

¹HBCDD and BDE-209 were determined in n=7 samples by GC-NCIMS following GC-ECD analysis.

Table S5. Recoveries of a OCPs and PBDE standard spike ($n = 1$) during analysis of the moose liver samples.

Analyte	%	Analyte	%	Analyte	%
PBDE 17	95	Hexachlorobutadiene	55	α -endosulfan	97
PBDE 28/33	101	1,2,4,5-TTCB	70	<i>cis</i> -Chlordane	72
PBDE 49	86	1,2,3,4-TTCB	68	Trans-nonachlor	67
PBDE 71	96	PECB	66	Dieldrin	79
PBDE 47	138	3,4,5-trichlorovertatrole	87	p,p'-DDE	67
PBDE 66	103	α -HCH	63	op-DDD	86
PBDE 100	116	β -HCH	78	Endrin	69
PBDE 99	133	HCB	71	b-Endosulfan	77
PBDE 85	107	3,4,5,6-tetrachlorovertatrole	76	<i>p,p'</i> -DDD	72
PBDE 154	100	Pentachloroanisole	68	<i>cis</i> -nonachlor	60
PBDE 153	104	γ -HCH (Indane)	67	α , p '-DDT	72
PBDE 138	85	Heptachlor	59	<i>p,p'</i> -DDT	68
PBDE 183	92	Aldrin	64	Methoxychlor	101
PBDE 190	92	Heptachlor Epoxide	79	Mirex	75
PBDE 209	71	Oxychlordane	81		
HBCDD	73	Octachlorostyrene	62		
BTBPE	78	trans-chlordane	62		
DBDPE	117	α , p '-DDE	72		

Table S6. Arithmetic and geometric mean concentrations of OCP/OCOs and PCBs in moose liver from the Dehcho and South Slave regions of the Northwest Territories (ng/g wet weight and lipid weight; n = 7 for each region). “<” values are instrument detection limits where arithmetic means were <0.001-0.002 ng/g ww.

	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	South	South	South	South	South	South	South	South	Slave
	Arith Mean	GM	min	max	Arith Mean	GM	min	max	Arith Mean	GM	min	GM	min	max	Slave	
Analyte	ng/ g ww	ng/ g ww	ng/ g ww	ng/ g ww	ng/g lw	ng/g lw	ng/g lw	ng/g lw	ng/ g ww	ng/ g ww	ng/ g ww	ng/g lw	ng/g lw	ng/g lw	ng/g lw	
% lipid	6.3	6.2	5.3	7.2					5.7	5.7	5.0	6.5				
Hexachlorobutadiene	0.011	0.010	0.007	0.014	0.17	0.12	0.24	0.005	0.004	0.003	0.011	0.080	0.074	0.050	0.16	
1,2,4,5-Tetrachloro-benzene	0.25	0.24	0.17	0.30	4.35	4.22	2.98	5.50	0.16	0.14	0.08	0.35	2.68	2.52	1.55	
1,2,3,4-Tetrachloro-benzene	0.28	0.27	0.18	0.35	4.85	4.68	3.16	5.87	0.16	0.15	0.09	0.36	2.79	2.64	1.86	
Pentachlorobenzene	0.11	0.037	0.008	0.32	1.85	0.59	0.11	5.91	0.17	0.16	0.10	0.45	3.00	2.74	1.75	
Hexachlorobenzene	0.17	0.17	0.10	0.24	2.85	2.67	1.45	4.33	0.34	0.26	0.11	0.93	6.02	4.64	1.62	
3,4,5,6-Tetrachloro-teratrole	<0.002				<0.02				<0.002	<0.002			<0.02			
Pentachloroanisole	0.040	0.017	0.005	0.15	0.70	0.28	0.07	2.66	0.029	0.023	0.004	0.07	0.52	0.41	0.08	
α -HCH	0.041	0.024	0.006	0.11	0.70	0.38	0.09	1.95	0.056	0.052	0.031	0.10	0.99	0.91	0.49	
β -HCH	0.10	0.097	0.059	0.17	1.66	1.56	1.11	2.93	0.079	0.072	0.040	0.17	1.38	1.27	0.71	
γ -HCH	0.056	0.035	0.005	0.14	0.95	0.57	0.07	2.48	0.063	0.058	0.032	0.13	1.11	1.02	0.56	
Pentachloronitrobenzene	<0.002				<0.02				<0.002	<0.002			<0.02			
Heptachlor	<0.002				<0.02				<0.002	<0.002			<0.02			
Aldrin	<0.002				<0.02				<0.002	<0.002			<0.02			
Dacthal	<0.002				<0.02				<0.002	<0.002			<0.02			
octachlorostyrene	<0.002				<0.02				<0.002	<0.002			<0.02			
Heptachlor Epoxide	0.044	0.033	0.010	0.080	0.716	0.53	0.16	1.47	0.046	0.037	0.017	0.15	0.78	0.65	0.27	
															2.23	

Table S6 continued . . .

Table S6 continued

	Dehcho	South Slave													
Analyte	ng/g ww	ng/g ww	ng/g ww	ng/g lw	ng/g ww	ng/g ww	ng/g lw								
Oxychlordane	0.033	0.020	0.004	0.10	0.510	0.32	0.064	1.48	0.039	0.025	0.009	0.14	0.66	0.44	0.15
trans-chlordane	<0.002				<0.02				<0.002	<0.002			<0.02		
cis-chlordane	0.017	0.011	<0.002	0.038	0.298	0.174	0.014	0.697	0.026	0.022	0.010	0.065	0.45	0.383	0.148
trans-nonachlor	0.031	0.016	0.004	0.058	<0.02				<0.002				<0.02		
Dieldrin	<0.002				<0.02				<0.002				<0.02		
cis-nonachlor	<0.002				<0.02				<0.002				<0.02		
Endrin	<0.002				<0.02				<0.002				<0.02		
α-Endosulfan	0.013	0.010	0.004	0.024	0.204	0.16	0.053	0.35	0.006	0.004	<0.002	0.025	0.11	0.066	0.021
β-Endosulfan	0.007	0.006	0.003	0.015	0.109	0.092	0.036	0.22	0.005	0.005	0.005	0.089	0.088	0.077	0.10
Endosulfan sulfate	0.026	0.015	0.003	0.092	0.434	0.24	0.059	1.73	0.005	0.004	0.002	0.008	0.086	0.079	0.034
<i>o,p'</i> -DDE	<0.002				<0.02				<0.02				<0.2		
<i>p,p'</i> -DDE	0.013	0.013	0.013	0.014	<0.02				<0.02				<0.2		
<i>o,p'</i> -DDD	<0.002				<0.02				<0.02				<0.2		
<i>p,p'</i> -DDD	0.017	0.017	0.017	0.017	<0.02				<0.02				<0.2		
<i>o,p'</i> -DDT	<0.002				<0.02				<0.02				<0.2		
<i>p,p'</i> -DDT	<0.002				<0.02				<0.02				<0.2		
Methoxychlor	<0.002				<0.02				<0.02				<0.2		
Mirex	0.004	0.002	<0.002	0.014	0.071	0.040	0.014	0.26	0.004	0.003	<0.002	0.013	0.073	0.049	0.012
Toxaphene	0.840	0.642	0.278	2.17	13.6	10.3	4.37	30.7	1.06	0.850	0.167	1.83	18.8	15.0	2.57
Hexachlorobornanes	0.026	0.008	<0.002	0.103	0.419	0.131	0.014	1.46	0.012	0.007	0.005	0.062	0.235	0.126	0.077

Table S6 continued . . .

Table S6 continued

	Dehcho	South	South	South	South	South	South	South	Slave							
Analyte	ng/g ww	Arith Mean	GM	min	max	Arith Mean	GM	min	max							
Heptachlorobornanes	0.366	0.047	0.003	1.50	5.83	0.753	0.039	21.3	0.123	0.039	0.005	0.494	2.077	0.682	0.077	7.72
Octachlorobornanes	0.086	0.039	<0.002	0.251	1.50	0.629	0.014	4.72	0.267	0.174	0.005	0.431	4.70	3.07	0.077	7.70
Nonachlorobornanes	0.318	0.311	0.225	0.447	5.10	4.99	3.54	6.72	0.481	0.260	0.005	1.26	8.50	4.60	0.077	22.6
Decachlorobornanes	0.048	0.043	0.015	0.085	0.770	0.683	0.211	1.183	0.186	0.104	0.010	0.525	3.43	1.83	0.208	10.5
P26	<0.02							<0.02					<0.2			
P50	<0.02							<0.02					<0.2			
P62	<0.02							<0.02					<0.2			
PCB-1	<0.002					<0.02			<0.002				<0.02			
PCB-3	<0.002				<0.02				<0.002				<0.02			
PCB4/10	0.005	0.005	0.004	0.009	0.086	0.083	0.051	0.13	0.028	0.008	0.005	0.19	0.502	0.14	0.077	3.39
PCB7/9	0.036	0.020	0.005	0.064	0.553	0.328	0.079	1.07	0.005	0.005	0.005	0.089	0.088	0.077	0.077	0.100
PCB6	0.006	0.006	0.004	0.008	0.092	0.090	0.060	0.13	0.005	0.005	0.005	0.089	0.088	0.077	0.077	0.100
PCB8/5	0.016	0.013	0.005	0.026	0.213	0.150	0.020	0.37	0.005	0.005	0.005	0.089	0.088	0.077	0.077	0.100
PCB12/13	<0.002				<0.02				<0.002				<0.02			
PCB15	0.117	0.011	<0.002	0.420	2.05	0.172	0.014	7.71	0.106	0.048	0.005	0.330	1.907	0.855	0.089	5.89
PCB19	<0.002				<0.02				<0.002	<0.002			<0.02			
PCB18	0.045	0.045	0.039	0.053	0.68	0.676	0.566	0.76	<0.002				<0.02			
PCB17	0.038	0.025	0.005	0.090	0.63	0.396	0.089	1.65	0.021	0.015	0.005	0.060	0.362	0.258	0.077	0.92
PCB27/24	0.003	0.002	<0.002	0.004	0.044	0.037	0.016	0.071	<0.002				<0.02			
PCB16/32	0.009	0.004	0.000	0.018	0.138	0.060	0.004	0.26	<0.002				<0.02			

Table S6 continued . . .

Table S6 continued

	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	South Slave						
Analyte	GM	min	max	Mean	GM	min	max	Mean	GM	min	max	Mean	GM	min	max
	ng/ g ww	ng/ g ww	ng/ g ww	ng/ g ww	ng/g lw	ng/g lw	ng/g lw	ng/g lw	ng/ g ww	ng/ g ww	ng/ g ww	ng/ g ww	ng/g lw	ng/g lw	ng/g lw
PCB26	<0.002			<0.02				<0.002	<0.002			<0.02			
PCB25	0.006	0.005	0.002	0.013	0.113	0.076	0.033	0.252	<0.002			<0.02			
PCB31/28	0.029	0.015	0.005	0.085	0.47	0.235	0.079	1.60	0.005	0.005	0.005	0.089	0.088	0.077	0.100
PCB20/33/21	0.043	0.014	0.004	0.24	0.74	0.217	0.057	4.26	0.068	0.015	0.005	0.435	1.061	0.259	0.078
PCB22	0.002	<0.002	0.003	0.027	0.020	0.010	0.044	<0.002				<0.02			6.69
PCB37	0.003	0.002	<0.002	0.005	0.046	0.032	0.014	0.092	0.018	0.007	0.005	0.110	0.351	0.130	0.077
PCB53	0.004	0.004	0.004	0.083	0.083	0.083	0.083	0.083	<0.002			<0.02			2.20
PCB45	0.011	0.010	0.008	0.013	0.18	0.170	0.120	0.24	<0.002			<0.02			
PCB46	0.006	0.005	0.003	0.009	0.10	0.084	0.044	0.16	<0.002			<0.02			
PCB73/52	0.079	0.020	<0.002	0.27	1.33	0.33	0.014	4.25	0.005	0.005	0.005	0.089	0.088	0.077	0.100
PCB43/49	<0.002			<0.02				<0.002	<0.002			<0.02			
PCB48/47/75	<0.002			<0.02				<0.002	<0.002			<0.02			
PCB44	<0.002			<0.02				<0.002	<0.002			<0.02			
PCB59/42	0.019	0.007	<0.002	0.036	0.298	0.118	0.016	0.67	<0.002	<0.002	<0.002	0.018	0.018	0.015	0.020
PCB71/41/68/64	<0.002			<0.02				<0.002	<0.002			<0.02			
PCB63	0.003	0.003	0.003	0.056	0.056	0.056	0.06	<0.002				<0.02			
PCB74/61	0.010	0.002	0.000	0.060	0.168	0.040	0.007	0.94	0.005	0.005	0.005	0.089	0.088	0.077	0.100
PCB70/76	<0.002			<0.02				<0.002	<0.002			<0.02			
PCB80/66	0.068	0.015	<0.002	0.13	1.054	0.24	0.016	2.49	<0.002	<0.002	<0.002	0.018	0.018	0.015	0.020
PCB56/60	0.041	0.008	<0.002	0.12	0.706	0.12	0.014	2.20	0.046	0.031	0.005	0.14	0.81	0.546	0.077

Table S6 continued . . .

Table S6 continued

Analyte	ng/g ww	ng/g ww	g ww	ng/g lw	g lw										
PCB81	<0.002	0.002	<0.002	0.005	0.037	0.026	0.014	0.092	0.005	0.005	0.005	<0.002	<0.002	<0.02	<0.02
PCB95/93	0.002	0.002	<0.002	0.005	<0.02	<0.02	<0.02	<0.02	<0.002	<0.002	<0.005	0.089	0.088	0.077	0.100
PCB91	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02	<0.02	<0.002	<0.002	<0.005	<0.002	<0.02	<0.02	<0.02
PCB92	<0.002	<0.002	<0.002	0.003	<0.002	0.100	0.286	0.040	0.014	1.83	0.005	0.005	0.005	0.089	0.088
PCB84/90/101/89	0.018	0.003	<0.002	0.002	<0.02	<0.02	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002
PCB99	<0.002	<0.002	<0.002	0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002
PCB83/108	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002
PCB97	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002
PCB86/111/125/117/87/ 116/115	<0.002	<0.002	0.008	0.075	0.452	0.122	0.014	1.38	0.028	0.025	0.015	0.040	0.45	0.23	0.71
PCB120/85	0.026	0.008	<0.002	0.075	0.452	0.122	0.014	1.38	0.028	0.025	0.015	0.040	0.45	0.23	0.71
PCB110	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002
PCB82	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002
PCB107/109	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002
PCB123	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002
PCB118/106	0.006	0.006	0.006	0.006	0.087	0.087	0.087	0.087	0.087	0.087	0.002	<0.002	<0.002	<0.002	<0.002
PCB114	0.000	0.000	0.000	0.000	0.003	0.003	0.003	0.003	0.003	0.003	<0.002	<0.002	<0.002	<0.002	<0.002
PCB105/127	0.002	<0.002	<0.002	0.006	0.031	0.021	0.014	0.12	<0.002	<0.002	<0.002	<0.002	<0.018	0.018	0.020
PCB126	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002
PCB151	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002	<0.002

Table S6 continued . . .

Table S6 continued

	Dehcho	South Slave													
Analyte	ng/g ww	ng/g ww	ng/g lw	ng/g ww	ng/g ww	ng/g ww	ng/g lw	ng/g lw	ng/g lw	ng/g lw					
PCB135/144	0.008	0.002	<0.002	0.051	0.149	0.028	0.014	0.95	<0.002	<0.002	<0.002	<0.002	0.018	0.015	0.020
PCB139/149	<0.002			<0.02					<0.002	<0.002			<0.02		
PCB131/165/142/146	0.003	0.002	<0.002	0.005	0.051	0.037	0.014	0.092	0.006	0.005	0.010	0.101	0.096	0.077	0.200
PCB153	0.003	0.002	<0.002	0.005	0.055	0.040	0.014	0.092	0.014	0.007	0.005	0.080	0.26	0.125	0.077
PCB132/168	<0.002			<0.02					<0.002	<0.002			<0.02		
PCB141	0.003	0.003	0.003	0.003	0.048	0.048	0.048	0.048	<0.002	<0.002			<0.02		
PCB137	0.003	0.002	<0.002	0.005	0.055	0.040	0.014	0.092	0.006	0.005	0.010	0.10	0.096	0.077	0.179
PCB163/164/138	<0.002			<0.02					<0.002	<0.002			<0.02		
PCB158/160	<0.002	<0.002	<0.002	<0.002	0.016	0.016	0.016	0.016	<0.002	<0.002			<0.02		
PCB129	0.003	0.003	0.003	0.003	0.058	0.058	0.058	0.058	<0.002	<0.002			<0.02		
PCB159	0.005	0.003	<0.002	0.012	0.076	0.049	0.016	0.22	<0.002	<0.002	<0.002	<0.002	0.018	0.015	0.020
PCB128/167	0.005	0.005	0.005	0.005	0.085	0.085	0.085	0.085	<0.002	<0.002	<0.002	<0.002	<0.02		
PCB156	0.002	0.002	<0.002	0.007	0.038	0.025	0.008	0.14	<0.002	<0.002	<0.002	<0.002	0.018	0.015	0.020
PCB157	<0.002			<0.02					<0.002	<0.002			<0.02		
PCB169	<0.002			<0.02					<0.002	<0.002			<0.02		
PCB182/187	0.003	0.002	<0.002	0.011	0.045	0.025	0.014	0.20	<0.002	<0.002	<0.002	<0.002	0.018	0.015	0.020
PCB183	0.003	0.002	<0.002	0.014	0.059	0.029	0.014	0.26	<0.002	<0.002	<0.002	<0.002	0.018	0.015	0.020
PCB174/181	<0.002			<0.02					<0.002	<0.002			<0.02		
PCB177	<0.002			<0.02					<0.002	<0.002			<0.02		
PCB171	<0.002			<0.02					<0.002	<0.002			<0.02		

Table S6 continued . . .

Table S6 continued

	Dehcho	South Slave											
Analyte	ng/g ww	ng/g ww	ng/g lw	ng/g ww	ng/g ww	ng/g lw	ng/g lw	ng/g lw	ng/g lw				
PCB172/192	<0.002	<0.002	<0.002	0.016	0.014	0.020	<0.002	<0.002	<0.002	<0.002	0.018	0.015	0.020
PCB180	0.003	0.002	<0.002	0.007	0.041	0.027	0.014	0.11	<0.002	<0.002	0.018	0.015	0.020
PCB193	0.009	0.002	<0.002	0.058	0.17	0.034	0.014	1.08	<0.002	<0.002	0.018	0.015	0.020
PCB191	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.011	<0.011	<0.002	<0.002	<0.02	<0.02	<0.02
PCB170/190	<0.002	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02	<0.002	<0.002	<0.02	<0.02	<0.02
PCB-202	<0.002	<0.002	<0.02	<0.02	<0.02	<0.030	0.016	0.090	<0.002	<0.002	<0.002	<0.02	<0.02
PCB199	<0.002	<0.002	<0.002	0.005	0.036	0.021	0.016	0.033	<0.002	<0.002	<0.002	<0.02	<0.02
PCB196/203	0.002	0.002	<0.002	0.002	0.022	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02
PCB195	<0.002	<0.002	<0.002	0.002	0.022	0.021	0.016	0.033	<0.002	<0.002	<0.002	<0.02	<0.02
PCB194	<0.002	<0.002	<0.002	0.002	0.022	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02
PCB205	<0.002	<0.002	<0.002	0.002	0.022	<0.02	<0.02	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02
PCB208	0.005	0.002	<0.002	0.032	0.099	0.027	0.009	0.593	<0.002	<0.002	0.018	0.015	0.020
PCB207	<0.002	<0.002	<0.002	0.032	0.099	0.027	0.009	0.593	<0.002	<0.002	<0.02	<0.02	<0.02
PCB206	<0.002	<0.002	<0.002	<0.002	<0.02	<0.002	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02
PCB209	<0.002	<0.002	<0.002	<0.002	<0.02	<0.002	<0.002	<0.002	<0.002	<0.002	<0.02	<0.02	<0.02

Table S7. Arithmetic and geometric mean concentrations of brominated and chlorinated flame retardants (ng/g wet wt and ng/g ww, $n = 7$). “<” values are instrument detection limits where arithmetic means were <0.001-0.002 ng/g ww. “na” indicates not available.

	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	Dehcho	South Slave					
	Arith Mean	GM	min	max	Arith Mean	GM	min	max	Arith Mean	GM	min	max	max
Analyte	ng/g ww	ng/g ww	ng/g ww	ng/g ww	ng/g lw	ng/g lw	ng/g lw	ng/g lw	ng/g lw	ng/g lw	ng/g lw	ng/g lw	ng/g lw
BDE 17	<0.002	0.001	<0.002	0.002	0.011	0.011	<0.007	0.022	<0.002	0.001	<0.002	0.009	<0.008
BDE 28/33	0.003	0.001	0.001	0.010	0.046	0.023	0.007	0.14	0.001	0.001	0.009	0.008	0.010
BDE 47	0.127	0.037	0.005	0.513	2.17	0.585	0.069	9.40	0.057	0.048	0.023	0.133	0.998
BDE 49	0.004	0.002	0.001	0.012	0.07	0.03	0.007	0.22	0.001	0.001	0.001	0.009	0.008
BDE 66	0.002	0.001	0.000	0.005	0.032	0.018	0.004	0.076	0.001	0.001	0.001	0.009	0.008
BDE 85	0.006	0.001	0.001	0.038	0.111	0.021	0.007	0.689	0.004	0.001	0.015	0.063	0.020
BDE 99	0.148	0.036	0.005	0.668	2.56	0.581	0.069	12.26	0.081	0.054	0.019	0.266	1.41
BDE 100	0.026	0.008	0.001	0.134	0.466	0.129	0.007	2.47	0.018	0.012	0.004	0.053	0.304
BDE 138	<0.002				0.111	0.021	0.007	0.69	<0.002			0.063	0.020
BDE 153	0.014	0.007	0.001	0.061	0.250	0.106	0.008	1.12	0.008	0.003	0.001	0.025	0.135
BDE 154	0.010	0.004	0.001	0.048	0.181	0.063	0.008	0.883	0.005	0.001	0.001	0.020	0.086
BDE 183	<0.002			<0.01					<0.002			<0.01	
BDE 190	<0.002			<0.01					<0.002			<0.01	
BDE 209	0.15	0.077	0.004	0.543	2.26	1.24	0.077	7.69	na	na	na	na	
TBP-AE	<0.002												3.57
TBP-DBPE	<0.002												0.389
BEHTBP	<0.002												0.320

Table S7 continued . . .

Table S7 continued

	Dehcho	South Slave													
Arith Mean	GM	min	max	Mean	GM	min	max	Mean	GM	min	max	Mean	GM	min	max
Analyte	ng/g ww	ng/g ww	ng/g ww	ng/g lw	ng/g ww	ng/g ww	ng/g ww	ng/g lw	ng/g lw	ng/g lw	ng/g lw				
BTBPE	<0.002			<0.01				<0.002				<0.01			<0.01
DPTE	<0.002			<0.01				<0.002				<0.01			<0.01
EHTeBB	<0.002			<0.01				<0.002				<0.01			<0.01
HBB	<0.002			<0.01				<0.002				<0.01			<0.01
OBIND	<0.002			<0.01				<0.002				<0.01			<0.01
PBBA	<0.002			<0.01				<0.002				<0.01			<0.01
PBBe	<0.002			<0.01				<0.002				<0.01			<0.01
PBEB	<0.002			<0.01				<0.002				<0.01			<0.01
PBTo	<0.002			<0.01				<0.002				<0.01			<0.01
pTBX	<0.002			<0.01				<0.002				<0.01			<0.01
syn-DP	<0.002			<0.01				<0.002				<0.01			<0.01
anti-DP	<0.002			<0.01				<0.002				<0.01			<0.01
TBCT	<0.002			<0.01				<0.002				<0.01			<0.01
HBCD	<0.002			<0.01				<0.002				<0.01			<0.01
BB-101	<0.002			<0.01				<0.002				<0.01			<0.01

Table S8. Arithmetic and geometric mean concentrations of individual PFASs in moose liver samples (ng/g ww).

Analyte	Dechcho			Dechcho			South Slave			South Slave		
	ng/g ww	ng/g ww	GM	ng/g ww	ng/g ww	max	ng/g ww	ng/g ww	AM	GM	ng/g ww	ng/g ww
PFBA	<0.16						<0.16					
PFPeA	<1.2						<1.2					
PFHxA	<0.15						<0.15					
PFHpA	<0.005						<0.005					
PFOA	0.069	0.067	0.056	0.113	0.090		0.085	0.050				
PFNA	0.255	0.243	0.174	0.448	0.188		0.171	0.080				
PFDA	0.240	0.204	0.078	0.601	0.159		0.149	0.069				
PFUnA	0.167	0.152	0.096	0.275	0.094		0.090	0.058				
PFDoA	0.008	0.005	0.003	0.020	0.017		0.016	0.008				
PFTriA	0.011	0.006	0.003	0.032	0.023		0.021	0.010				
PFTA	0.004	0.003	<0.003	0.015	0.012		0.008	0.003				
PFHxDAA	0.004	0.003	0.003	0.010	0.049		0.032	0.010				
PFBS	0.055	0.052	0.026	0.087	0.029		0.018	0.003				
PFHxS	0.063	0.057	0.022	0.106	0.012		0.007	0.003				
PFHpS	<0.005						<0.005					
PFOS	0.437	0.377	0.210	0.993	0.252		0.244	0.174				
PFOS_Linear	0.189	0.189	0.173	0.202	0.183		0.172	0.088				
PFDS	0.009	0.005	0.003	0.034	<0.005							
PFOSA	0.008	0.004	0.003	0.030	0.012		0.008	0.003				

Table S9A. Pearson correlation matrix for major halogenated organic contaminants in moose liver. Correlation coefficients in bold are significant at $P < 0.05$.

	Age	% lipid	Log Σ PFCA	Log Σ PFSA	Log Σ_{13} PBDE	Log Σ PCB	Log Σ Endo-sulfan	Log Toxaphene	Log Σ DDT	Log Σ CHL	Log Σ HCH	Log Σ CBZ	Log Σ Mono_Di_CB	Log Σ Tri-CB	Log Σ Tetra-CB	Log Σ penta-CB	Log Σ hexa-CB
Age	1																
% lipid	0.266	1															
Log Σ PFCA	0.248	0.406	1														
Log Σ PFSA	0.076	0.590	0.699	1													
Log Σ_{13} PBDE	0.187	0.341	0.018	0.138	1												
Log Σ PCB	0.285	0.042	-0.001	-0.033	0.460	1											
Log Σ Endo-sulfan	0.006	0.341	0.328	0.618	0.005	0.309	1										
Log Toxaphene	-0.288	-0.152	-0.498	-0.505	-0.201	-0.046	-0.054	1									
Log Σ DDT	-0.368	0.42	-0.038	0.418	-0.086	0.103	0.623	0.268	1								
Log Σ CHL	0.107	0.103	0.05	-0.096	0.602	0.358	-0.423	-0.314	-0.505	1							
Log Σ HCH	0.149	-0.240	-0.251	-0.405	-0.011	-0.294	-0.79	-0.195	-0.698	0.514	1						
Log Σ CBZ	0.38	-0.496	-0.163	-0.555	0.032	0.052	-0.622	-0.154	-0.821	0.457	0.75	1					
Log Σ Mono_Di_CB	0.227	-0.556	-0.371	-0.829	-0.324	0.147	-0.514	0.212	-0.559	0.089	0.488	0.671	1				
Log Σ Tri-CB	-0.028	-0.481	-0.118	-0.583	-0.325	0.036	-0.643	-0.038	-0.502	0.321	0.594	0.592	0.781	1			
Log Σ Tetra-CB	0.485	-0.319	-0.426	-0.605	0.154	0.613	-0.285	0.091	-0.267	0.201	0.233	0.546	0.709	0.459	1		
Log Σ penta-CB	0.518	-0.384	-0.018	-0.339	-0.303	0.223	-0.202	-0.229	-0.426	0.059	0.344	0.628	0.681	0.664	0.667	1	
Log Σ hexa-CB	-0.294	0.421	0.365	0.576	-0.196	-0.081	0.555	0.277	0.686	-0.309	-0.659	-0.717	-0.662	-0.453	-0.588	-0.416	

Table S9B. Pearson correlation matrix for selected individual halogenated organic contaminants in moose liver. Correlation coefficients in bold are significant at $P < 0.05$.

	AGE	% lipid	Log HCBD	Log 1245-TeCBz	Log 1234-TeCBz	Log PeCBz	Log HCB	Log PCA	Log α HCH	Log β HCH	Log heptachlor epoxide	Log Oxychlordane	
AGE	1.000												
% lipid	0.266	1.000											
Log HCBD	0.035	0.569	1.000										
Log 1245_TeCBz	0.341	-0.513	-0.608	1.000									
Log 1234_TeCBz	0.347	-0.513	-0.599	1.000									
Log PeCBz	0.342	-0.457	-0.552	0.972	1.000								
Log HCB	0.103	-0.426	-0.510	0.467	0.458	0.476	1.000						
Log PCA	0.436	-0.298	-0.529	0.778	0.770	0.828	0.543	1.000					
Log α HCH	0.240	-0.587	-0.520	0.920	0.919	0.872	0.591	0.664	1.000				
Log β HCH	-0.178	0.408	0.416	-0.130	-0.128	-0.117	-0.341	-0.217	-0.170	1.000			
Log γ HCH	0.139	-0.351	-0.551	0.641	0.641	0.697	0.414	0.537	0.512	0.155	1.000		
Log heptachlor epoxide	0.329	0.021	-0.059	0.531	0.534	0.542	0.116	0.556	0.327	0.172	0.302	1.000	
Log Oxychlordane	-0.038	0.265	-0.027	0.114	0.112	0.212	0.335	0.308	0.146	0.282	0.364	0.306	
Log <i>cis</i> -chlordane	0.201	-0.594	-0.376	0.721	0.723	0.679	0.283	0.462	0.633	-0.232	0.383	0.586	
Log α -Endosulfan	0.124	0.369	0.327	-0.320	-0.314	-0.283	-0.457	0.023	-0.494	-0.025	-0.406	0.380	
Log endosulfan sulfate	0.072	0.432	0.635	-0.729	-0.726	-0.700	-0.499	-0.528	-0.650	-0.180	-0.693	-0.436	
Log mirex	-0.102	-0.464	0.036	0.186	0.197	0.127	-0.047	-0.274	0.321	0.134	0.182	-0.154	
Log BDE47	0.462	-0.225	-0.311	0.631	0.628	0.632	0.333	0.736	0.448	-0.068	0.425	0.830	
Log PFNA	0.148	0.452	0.548	-0.311	-0.306	-0.285	-0.498	-0.336	-0.437	0.415	-0.071	0.333	
Log PFDA	0.391	0.339	0.295	-0.211	-0.208	-0.264	-0.010	-0.234	-0.246	0.152	-0.100	0.296	
Log PFUNA	0.025	0.502	0.625	-0.677	-0.671	-0.694	-0.433	-0.665	-0.705	0.085	-0.617	-0.010	-0.282
Log PFBS	-0.005	0.245	0.589	-0.185	-0.170	-0.173	-0.475	-0.246	-0.201	0.743	-0.101	0.213	0.001
Log PFHXS	0.193	0.359	0.678	-0.408	-0.397	-0.336	-0.164	-0.149	-0.281	-0.160	-0.627	0.053	0.151
Log PFOS	-0.008	0.555	0.611	-0.705	-0.704	-0.715	-0.236	-0.590	-0.702	0.368	-0.381	-0.023	-0.087

	Log Oxychlordane	Log <i>cis</i> -chlordane	Log α Endosulfan	Log endosulfan sulfate	Log BDE47	Log PFNA	Log PFDA	Log PFUNA	Log PFBS	Log PFHXS
AGE										
% lipid										
Log HCBD										
Log 1245_TCBZ										
Log 1234_TCBZ										
Log PeCBZ										
Log HCB										
Log PCA										
Log α HCH										
Log β HCH										
Log γ HCH										
Log heptachlor epoxide										
Log Oxychlordane										
Log <i>cis</i> -chlordane	1									
Log α Endosulfan	−0.076									
Log endosulfan sulfate	−0.353	0.459	1							
Log minex	0.444	−0.471	−0.044							
Log BDE47	0.635	0.119	−0.531	−0.087						
Log PFNA	0.151	0.302	0.346	0.238	0.18					
Log PFDA	0.185	0.128	0.212	0.228	0.293	0.73				
Log PFUNA	−0.152	0.382	0.607	0.053	−0.161	0.695	0.717	1		
Log PFBS	−0.169	0.323	0.085	0.276	−0.044	0.38	0.182	0.237	1	
Log PFHXS	−0.267	0.482	0.552	−0.209	−0.085	0.126	0.119	0.419	0.272	1
Log PFOS	−0.317	0.313	0.487	−0.015	−0.185	0.658	0.759	0.842	0.369	0.257

Table S9C. Pearson correlation matrix for selected PCB congeners (co-eluters reported as single GC peaks) in moose liver. Correlation coefficients in bold are significant at $P < 0.05$.

	Age	%Lipid	Log CB 15	Log CB 31_28	Log CB 73_52	Log CB 56_60	Log CB 105_127	Log CB153	Log CB156	Log CB180
Age	1									
%Lipid	0.266	1								
Log CB15	-0.455	0.066	1							
Log CB31_28	0.388	-0.424	-0.309	1						
Log CB73_52	-0.372	0.351	0.155	-0.721	1					
Log CB56_60	0.395	-0.054	-0.285	0.249	0.213	1				
Log CB105_127	0.424	-0.532	-0.187	0.922	-0.770	0.243	1			
Log CB153	-0.239	-0.223	-0.003	-0.367	0.646	0.220	-0.311	1		
Log CB156	0.129	-0.572	0.250	0.326	-0.466	0.043	0.625	0.014	1	
Log CB180	-0.113	0.021	-0.025	-0.500	0.474	-0.123	-0.492	0.796	-0.194	1
Log CB196_203	-0.063	0.139	-0.212	-0.538	0.438	-0.181	-0.558	0.612	-0.299	0.886

Table S10A. Comparison of mean concentrations (log ng/g wet wt) of major halogenated organics in moose liver from the Dehcho and South Slave (SSR) regions using the Students *t*-test assuming separate variance; significant values are bolded ($P < 0.05$).

Variable	Region	n	Mean	Stand Dev	Mean Difference	Lower 95% Limit	Upper 95% Limit	t	df	p-Value
Log Σ PFCA	Dehcho	7	-0.152	0.162	0.057	-0.105	0.218	0.783	10,129	0.452
	SSR	7	-0.209	0.102						
Log Σ PFSA	Dehcho	7	-0.284	0.189	0.245	0.061	0.429	3.0	9,382	0.014
	SSR	7	-0.529	0.105						
Log Σ PBDE	Dehcho	7	-0.589	0.516	0.354	-0.175	0.884	1.474	10,923	0.169
	SSR	7	-0.943	0.373						
Log Σ PCB	Dehcho	7	-0.217	0.177	0.24	0.016	0.465	2.341	11,742	0.038
	SSR	7	-0.458	0.206						
Log Σ Endosulfan	Dehcho	7	-1.478	0.373	0.361	-0.002	0.724	2.238	9,396	0.051
	SSR	7	-1.839	0.208						
Log Toxaphene	Dehcho	7	-0.192	0.329	-0.190	-0.533	0.153	-1.215	11,197	0.249
	SSR	7	-0.002	0.250						
Log Σ CHL	Dehcho	7	-1.052	0.296	0.033	-0.281	0.346	0.23	11	0.823
	SSR	7	-1.085	0.237						
Log Σ HCH	Dehcho	7	-0.859	0.143	-0.123	-0.277	0.03	-1.76	11,553	0.105
	SSR	7	-0.736	0.118						
Log Σ CBZ	Dehcho	7	-0.479	0.443	-0.351	-0.767	0.066	-1.954	7,722	0.088
	SSR	7	-0.128	0.17						
Log Σ Mono_Di-CB	Dehcho	7	-1.503	1.086	-1.081	-2.118	-0.04	-2.375	8,616	0.043
	SSR	7	-0.422	0.520						
Log Σ Tri-CB	Dehcho	7	-1.844	0.830	-0.608	-1.402	0.187	-1.739	8,735	0.117
	SSR	7	-1.236	0.408						
Log Σ Tetra-CB	Dehcho	7	-1.301	0.690	0.009	-0.639	0.658	0.033	7,712	0.975
	SSR	7	-1.31	0.264						
Log Σ hexa-CB	Dehcho	7	-1.349	0.621	0.342	-0.251	0.936	1.314	8,635	0.223
	SSR	7	-1.692	0.299						

Table S10B. Comparison of mean concentrations ($\log \text{ng/g}$ wet wt) of major individual halogenated organic contaminants in moose liver from the Dehcho and South Slave (SSR) regions using the Students *t*-test assuming separate variance; significant values are bolded ($P < 0.05$).

Variable	Region	n	Mean	Stand Dev	Mean Difference	Lower 95% CI	Upper 95% CI	t	df	p-Value
Log HCBD	Dehcho	7	-1.982	0.13	0.406	0.254	0.558	5.81	12	0.001
	SSR	7	-2.388	0.131						
Log 1245-TeCBz	JMR	7	-1.804	1.119	-0.932	-1.968	0.103	-2.184	6.224	0.07
	SSR	7	-0.872	0.153						
Log 1234-TeCBz	JMR	7	-1.785	1.144	-0.932	-1.99	0.126	-2.139	6.181	0.075
	SSR	7	-0.853	0.14						
Log PeCBz	Dehcho	7	-1.436	0.725	-0.599	-1.271	0.073	-2.14	7	0.072
	SSR	7	-0.837	0.16						
Log HCB	Dehcho	7	-0.779	0.135	-0.229	-0.505	0.046	-1.90	8	0.092
	SSR	7	-0.549	0.29						
Log PCA	Dehcho	7	-2.03	0.67	-0.397	-1.051	0.256	-1.36	10	0.204
	SSR	7	-1.633	0.381						
Log α -HCH	Dehcho	7	-1.623	0.513	-0.341	-0.818	0.137	-1.683	7.07	0.118
	SSR	7	-1.282	0.154						
Log β -HCH	Dehcho	7	-1.109	0.081	0.044	-0.108	0.197	0.657	8.92	0.523
	SSR	7	-1.153	0.159						
Log γ -HCH	Dehcho	7	-1.684	0.313	-0.44	-0.741	-0.138	-3.307	8.93	0.006
	SSR	7	-1.244	0.116						
Log heptachlor-epoxide	Dehcho	7	-1.480	0.387	-0.002	-0.385	0.38	-0.014	9.923	0.989
	SSR	7	-1.477	0.236						
Log oxychlordane	Dehcho	7	-1.7	0.508	-0.040	-0.577	0.497	-0.16	11	0.873
	SSR	7	-1.66	0.401						
Log <i>cis</i> -chlordane	Dehcho	7	-1.965	0.552	-0.280	-0.803	0.243	-1.23	8	0.254
	SSR	7	-1.686	0.241						

Table S10B continued . . .

Table S10B continued

Variable	Region	n	Mean	Stand Dev	Mean Difference	Lower 95% CI	Upper 95% CI	t	df	p-Value
Log α -endosulfan	Dehcho	7	-2.091	0.424	0.353	-0.147	0.854	1.54	12	0.150
Log endosulfan sulfate	SSR	7	-2.445	0.435						
Log endosulfan sulfate	Dehcho	7	-1.825	0.479	0.521	0.062	0.980	2.58	9	0.030
Log mirex	SSR	7	-2.346	0.236						
Log mirex	Dehcho	7	-2.604	0.437	-0.031	-0.536	0.474	-0.135	12	0.895
Log BDE47	SSR	7	-2.573	0.430						
Log BDE47	Dehcho	7	-1.437	0.851	-0.097	-0.890	0.697	-0.29	7	0.783
Log PFNA	SSR	7	-1.341	0.277						
Log PFNA	Dehcho	7	-0.615	0.142	0.153	-0.057	0.363	1.61	11	0.137
Log PFDA	SSR	7	-0.768	0.207						
Log PFDA	Dehcho	7	-0.69	0.26	0.138	-0.126	0.402	1.15	11	0.274
Log PFUNA	SSR	7	-0.827	0.179						
Log PFUNA	Dehcho	7	-0.817	0.2	0.226	0.028	0.425	2.54	10	0.029
Log PFOS	SSR	7	-1.044	0.126						
Log PFOS	Dehcho	7	-0.423	0.243	0.189	-0.043	0.421	1.86	9	0.098
Log PFBS	SSR	7	-0.612	0.117						
Log PFBS	Dehcho	7	-1.28	0.154	0.467	-0.078	1.012	2.035	6.827	0.082
Log PFHxS	SSR	7	-1.747	0.587						
Log PFHxS	Dehcho	7	-1.246	0.226	0.94	0.48	1.4	4.663	8.504	0.001
Log PCB15	SSR	7	-2.186	0.483						
Log PCB15	Dehcho	7	-1.969	1.29	-0.645	-1.903	0.614	-1.15	10	0.278
Log PCB17	SSR	7	-1.325	0.733						
Log PCB17	Dehcho	7	-1.607	0.454	0.199	-0.259	0.658	0.96	11	0.358
Log PCB17	SSR	7	-1.806	0.307						

Table S10B continued . . .

Table S10B continued

Variable	Region	n	Mean	Stand Dev	Mean Difference	Lower 95% CI	Upper 95% CI	t	df	p-Value
Log CB20_33_21	Dehcho	7	-1.869	0.616	0.064	-0.685	0.812	0.19	12	0.855
	SSR	7	-1.932	0.668						
Log PCB56_60	Dehcho	7	-2.109	1.01	-0.675	-1.619	0.268	-1.68	7	0.135
	SSR	7	-1.433	0.342						
Log PCB153	Dehcho	7	-2.607	0.368	-0.478	-0.962	0.006	-2.16	11	0.053
	SSR	7	-2.129	0.455						