

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD060525\
 Data File : PD088817.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 05 Jun 2025 17:33
 Operator : AR\AJ
 Sample : PIBLK174
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PIBLK527

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 06 11:17:13 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD060525CLP.M
 Quant Title : GC Extractables
 QLast Update : Fri Jun 06 11:09:32 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.549	2.880	59114689	354.3E6	23.095	23.137
2)	SA Decachlor...	9.072	8.073	137.6E6	693.4E6	35.524	35.337
Target Compounds							
3)	MA gamma-BHC...	4.340	0.000	384866	0	<MDL	N.D. #
6)	B beta-BHC	0.000	3.990f	0	18633645	N.D.	1.828
7)	B delta-BHC	4.775	0.000	2700370	0	0.569	N.D. #
8)	B Heptachlo...	5.646f	0.000	193435	0	<MDL	N.D. #
9)	A Endosulfan I	6.079	0.000	11722	0	<MDL	N.D. #
10)	B trans-Chl...	5.996f	0.000	2147167	0	0.505	N.D. #
11)	B cis-Chlor...	5.996	0.000	2147167	0	0.493	N.D. #
12)	B 4,4'-DDE	6.215	0.000	364638	0	<MDL	N.D. #
13)	MA Dieldrin	6.404f	0.000	859901	0	0.196	N.D. #
14)	MA Endrin	6.590	0.000	580073	0	0.157	N.D. #
15)	B Endosulfa...	6.772	0.000	2643722	0	0.696	N.D. #
16)	A 4,4'-DDD	6.772f	0.000	2643722	0	0.846	N.D. #
22)	Toxaphene-1	6.215	0.000	364638	0	12.814	N.D. #
23)	Toxaphene-2	6.404f	0.000	859901	0	21.518	N.D. #

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

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