

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD110119\
 Data File : PD055733.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 31 Oct 2019 01:43
 Operator : SG\AJ
 Sample : PCHLORICV500
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 ICVPD110119CHLOR

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Oct 31 07:42:26 2019
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD110119.M
 Quant Title : GC Extractables
 QLast Update : Thu Oct 31 06:48:28 2019
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 µl
 Signal #1 Phase : ZB-MR2 Signal #2 Phase: ZB-MR1
 Signal #1 Info : 30M x 0.32mm x0.2 Signal #2 Info : 30M x 0.32mm x 0.50µm

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 15% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(Min)
1 SA	Tetrachloro-m-xylene	852.507	874.195 E3	-2.5	107	0.00
23	Chlordane-1	25.934	25.345 E3	2.3	96	0.00
24	Chlordane-2	31.802	31.236 E3	1.8	102	0.00
25	Chlordane-3	96.166	90.956 E3	5.4	96	0.00
26	Chlordane-4	84.097	79.422 E3	5.6	96	0.00
27	Chlordane-5	30.364	30.340 E3	0.1	99	0.00
28 SA	Decachlorobiphenyl	946.056	870.814 E3	8.0	97	0.00

Signal #2

1 SA	Tetrachloro-m-xylene	1.204	1.129 E6	6.2	98	0.00
23	Chlordane-1	46.406	45.630 E3	1.7	101	0.00
24	Chlordane-2	52.530	48.782 E3	7.1	98	0.00
25	Chlordane-3	218.980	212.205 E3	3.1	99	0.00
26	Chlordane-4	263.049	252.684 E3	3.9	99	0.00
27	Chlordane-5	46.619	45.847 E3	1.7	100	0.00
28 SA	Decachlorobiphenyl	1.392	1.266 E6	9.1	96	0.00

Evaluate Continuing Calibration Report - Not Found

2 MA	alpha-BHC	1.270	0.000 E6	100.0#	0#	-3.72#
3 MA	gamma-BHC (Lindane)	1.198	0.000 E6	100.0#	0#	-3.99#
4 MA	Heptachlor	1.065	0.000 E6	100.0#	0#	-4.28#
5 MB	Aldrin	1.094	0.000 E6	100.0#	0#	-4.52#
6 B	beta-BHC	504.056	0.000 E3	100.0#	0#	-4.24#
7 B	delta-BHC	1.152	0.000 E6	100.0#	0#	-4.44#
8 B	Heptachlor epoxide	1.029	0.000 E6	100.0#	0#	-4.95#
9 A	Endosulfan I	974.863	0.000 E3	100.0#	0#	-5.28#
10 B	gamma-Chlordane	1.067	0.000 E6	100.0#	0#	-5.18#
11 B	alpha-Chlordane	1.055	0.000 E6	100.0#	0#	-5.23#
12 B	4,4'-DDE	984.301	0.000 E3	100.0#	0#	-5.40#
13 MA	Dieldrin	1.061	0.000 E6	100.0#	0#	-5.52#
14 MA	Endrin	959.019	0.000 E3	100.0#	0#	-5.78#
15 B	Endosulfan II	969.226	0.000 E3	100.0#	0#	-6.05#
16 A	4,4'-DDD	803.984	0.000 E3	100.0#	0#	-5.91#
17 MA	4,4'-DDT	715.629	0.000 E3	100.0#	0#	-6.15#
18 B	Endrin aldehyde	835.661	0.000 E3	100.0#	0#	-6.22#
19 B	Endosulfan Sulfate	895.779	0.000 E3	100.0#	0#	-6.43#
20 A	Methoxychlor	391.837	0.000 E3	100.0#	0#	-6.70#
21 B	Endrin ketone	1.046	0.000 E6	100.0#	0#	-6.91#
22	Mirex	824.578	0.000 E3	100.0#	0#	-7.09#

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(Min)

Signal #2					
2 A alpha-BHC	1.762	0.000 E6	100.0#	0#	-4.35#
3 MA gamma-BHC (Lindane)	1.717	0.000 E6	100.0#	0#	-4.63#
4 MA Heptachlor	1.733	0.000 E6	100.0#	0#	-5.14#
5 MB Aldrin	1.674	0.000 E6	100.0#	0#	-5.44#
6 B beta-BHC	794.843	0.000 E3	100.0#	0#	-4.80#
7 B delta-BHC	1.603	0.000 E6	100.0#	0#	-5.02#
8 B Heptachlor epoxide	1.615	0.000 E6	100.0#	0#	-5.83#
9 A Endosulfan I	1.571	0.000 E6	100.0#	0#	-6.18#
10 B gamma-Chlordane	1.724	0.000 E6	100.0#	0#	-6.06#
11 B alpha-Chlordane	1.718	0.000 E6	100.0#	0#	-6.13#
12 B 4,4'-DDE	1.599	0.000 E6	100.0#	0#	-6.29#
13 MA Dieldrin	1.690	0.000 E6	100.0#	0#	-6.44#
14 MA Endrin	1.442	0.000 E6	100.0#	0#	-6.65#
15 B Endosulfan II	1.553	0.000 E6	100.0#	0#	-6.86#
16 A 4,4'-DDD	1.366	0.000 E6	100.0#	0#	-6.77#
17 MA 4,4'-DDT	1.335	0.000 E6	100.0#	0#	-7.07#
18 B Endrin aldehyde	1.341	0.000 E6	100.0#	0#	-6.98#
19 B Endosulfan Sulfate	1.455	0.000 E6	100.0#	0#	-7.21#
20 A Methoxychlor	799.943	0.000 E3	100.0#	0#	-7.53#
21 B Endrin ketone	1.659	0.000 E6	100.0#	0#	-7.68#
22 Mirex	1.210	0.000 E6	100.0#	0#	-8.13#

(#) = Out of Range

SPCC's out = 0 CCC's out = 0