Evaluate Continuing Calibration Report

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 06:45:42 2019
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD110119.M
Quant Title : GC Extractables
QLast Update : Thu Oct 31 06:44:18 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	923.500	876.927 E3	5.0	98	0.00
4	MA	Heptachlor	0.000	382587.634	0.0	0#	0.00
5	MB	Aldrin	0.000	33122.314	0.0	0#	0.01
7	B	delta-BHC	0.000	33821.193	0.0	0#	0.00
9	A	Endosulfan I	0.000	75678.678	0.0	0#	-0.01
10	B	gamma-Chlordane	0.000	935163.476	0.0	0#	0.00
11	B	alpha-Chlordane	0.000	823151.998	0.0	0#	0.00
14	MA	Endrin	0.000	63942.853	0.0	0#	0.00
15	B	Endosulfan II	0.000	303404.698	0.0	0#	0.00
18	B	Endrin aldehyde	0.000	16574.949	0.0	0#	0.00
23		Chlordane-1	25.934	26.361 E3	-1.6	100	0.00
24		Chlordane-2	31.802	31.855 E3	-0.2	104	0.00
25		Chlordane-3	96.166	93.516 E3	2.8	99	0.00
26		Chlordane-4	84.097	82.315 E3	2.1	99	0.00
27		Chlordane-5	30.364	30.340 E3	0.1	99	0.00
28	SA	Decachlorobiphenyl	950.478	870.024 E3	8.5	96	0.00

Signal #2

1	SA	Tetrachloro-m-xylene	1.196	1.130 E6	5.5	98	0.00
2	A	alpha-BHC	0.000	35634.718	0.0	0#	0.00
4	MA	Heptachlor	0.000	702138.441	0.0	0#	0.00
5	MB	Aldrin	0.000	490837.379	0.0	0#	-0.03
6	B	beta-BHC	0.000	34059.605	0.0	0#	0.03
8	B	Heptachlor epoxide	0.000	431882.809	0.0	0#	-0.01
10	B	gamma-Chlordane	0.000	2158412.030	0.0	0#	0.00
11	B	alpha-Chlordane	0.000	2599187.636	0.0	0#	0.00
13	MA	Dieldrin	0.000	124033.064	0.0	0#	0.00
14	MA	Endrin	0.000	77908.325	0.0	0#	0.02
15	B	Endosulfan II	0.000	52129.352	0.0	0#	0.01
16	A	4,4'-DDD	0.000	148945.261	0.0	0#	0.02
23		Chlordane-1	46.406	45.749 E3	1.4	101	0.00
24		Chlordane-2	52.530	49.084 E3	6.6	98	0.00
25		Chlordane-3	218.980	215.841 E3	1.4	101	0.00
26		Chlordane-4	263.049	259.919 E3	1.2	102	0.00
27		Chlordane-5	46.619	51.176 E3	-9.8	111	0.00
28	SA	Decachlorobiphenyl	1.390	1.264 E6	9.1	96	0.00

Evaluate Continuing Calibration Report - Not Founds

2	MA	alpha-BHC	0.000	0.000	0.0	0# -3.72#
3	MA	gamma-BHC (Lindane)	0.000	0.000	0.0	0# -3.99#

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 06:45:42 2019
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD110119.M
Quant Title : GC Extractables
QLast Update : Thu Oct 31 06:44:18 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)
6	B	beta-BHC	0.000	0.000	0.0	0#	-4.24#
8	B	Heptachlor epoxide	0.000	0.000	0.0	0#	-4.95#
12	B	4,4'-DDE	0.000	0.000	0.0	0#	-5.40#
13	MA	Dieldrin	0.000	0.000	0.0	0#	-5.52#
16	A	4,4'-DDD	0.000	0.000	0.0	0#	-5.91#
17	MA	4,4'-DDT	0.000	0.000	0.0	0#	-6.15#
19	B	Endosulfan Sulfate	0.000	0.000	0.0	0#	-6.43#
20	A	Methoxychlor	0.000	0.000	0.0	0#	-6.70#
21	B	Endrin ketone	0.000	0.000	0.0	0#	-6.91#
22		Mirex	0.000	0.000	0.0	0#	-7.09#

Signal #2

3	MA	gamma-BHC (Lindane)	0.000	0.000	0.0	0#	-4.63#
7	B	delta-BHC	0.000	0.000	0.0	0#	-5.02#
9	A	Endosulfan I	0.000	0.000	0.0	0#	-6.18#
12	B	4,4'-DDE	0.000	0.000	0.0	0#	-6.29#
17	MA	4,4'-DDT	0.000	0.000	0.0	0#	-7.07#
18	B	Endrin aldehyde	0.000	0.000	0.0	0#	-6.98#
19	B	Endosulfan Sulfate	0.000	0.000	0.0	0#	-7.21#
20	A	Methoxychlor	0.000	0.000	0.0	0#	-7.53#
21	B	Endrin ketone	0.000	0.000	0.0	0#	-7.68#
22		Mirex	0.000	0.000	0.0	0#	-8.13#

(#) = Out of Range SPCC's out = 0 CCC's out = 0