Evaluate Continuing Calibration Report

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL080521\
Data File : PL068717.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On   : 05 Aug 2021 15:11
Operator : AR\AJ
Sample   : PCHLORICV500
Misc   :
ALS Vial : 21   Sample Multiplier: 1

Instrument : ECD_L
ClientSampleId : ICVPL080521CHLOR

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Aug 06 03:07:29 2021
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL080521.M
Quant Title : GC Extractables
QLast Update : Fri Aug 06 03:02:43 2021
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj.   : 1 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 x0.5µm

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

		Compound	Amount	Calc.	%Dev	Area%	Dev(Min)
1	SA	Tetrachloro-m-xylene	50.000	54.670	-9.3	110	0.00
23		Chlordane-1	500.000	504.809	-1.0	100	0.00
24		Chlordane-2	500.000	489.169	2.2	99	0.00
25		Chlordane-3	500.000	498.067	0.4	99	0.00
26		Chlordane-4	500.000	496.497	0.7	99	0.00
27		Chlordane-5	500.000	485.085	3.0	99	0.00
28	SA	Decachlorobiphenyl	50.000	49.551	0.9	103	0.00

Signal #2

1	SA	Tetrachloro-m-xylene	50.000	52.690	-5.4	104	0.00
23		Chlordane-1	500.000	480.554	3.9	100	0.00
24		Chlordane-2	500.000	505.806	-1.2	99	0.00
25		Chlordane-3	500.000	497.606	0.5	98	0.00
26		Chlordane-4	500.000	496.794	0.6	98	0.00
27		Chlordane-5	500.000	498.829	0.2	97	0.00
28	SA	Decachlorobiphenyl	50.000	51.017	-2.0	102	0.00

Evaluate Continuing Calibration Report - Not Founds

2	A	alpha-BHC	50.000	0.000	100.0#	0	-5.35#
3	MA	gamma-BHC (Lindane)	50.000	0.000	100.0#	0	-5.77#
4	MA	Heptachlor	50.000	0.000	100.0#	0	-6.18#
5	MB	Aldrin	50.000	0.000	100.0#	0	-6.51#
6	B	beta-BHC	50.000	0.000	100.0#	0	-6.15#
7	B	delta-BHC	50.000	0.000	100.0#	0	-6.41#
8	B	Heptachlor epoxide	50.000	0.000	100.0#	0	-7.08#
9	A	Endosulfan I	50.000	0.000	100.0#	0	-7.49#
10	B	gamma-Chlordane	50.000	0.000	100.0#	0	-7.36#
11	B	alpha-Chlordane	50.000	0.000	100.0#	0	-7.43#
12	B	4,4'-DDE	50.000	0.000	100.0#	0	-7.65#
13	MA	Dieldrin	50.000	0.000	100.0#	0	-7.77#
14	MA	Endrin	50.000	0.000	100.0#	0	-8.06#
15	B	Endosulfan II	50.000	0.000	100.0#	0	-8.37#
16	A	4,4'-DDD	50.000	0.000	100.0#	0	-8.23#
17	MA	4,4'-DDT	50.000	0.000	100.0#	0	-8.49#
18	B	Endrin aldehyde	50.000	0.000	100.0#	0	-8.56#
19	B	Endosulfan Sulfate	50.000	0.000	100.0#	0	-8.79#
20	A	Methoxychlor	50.000	0.000	100.0#	0	-9.09#
21	B	Endrin ketone	50.000	0.000	100.0#	0	-9.31#
22		Mirex	50.000	0.000	100.0#	0	-9.50#

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Max. RRF Dev : 15%   Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(Min)
Signal #2							
2	A	alpha-BHC	50.000	0.000	100.0#	0	-6.15#
3	MA	gamma-BHC (Lindane)	50.000	0.000	100.0#	0	-6.55#
4	MA	Heptachlor	50.000	0.000	100.0#	0	-7.20#
5	MB	Aldrin	50.000	0.000	100.0#	0	-7.57#
6	B	beta-BHC	50.000	0.000	100.0#	0	-6.80#
7	B	delta-BHC	50.000	0.000	100.0#	0	-7.07#
8	B	Heptachlor epoxide	50.000	0.000	100.0#	0	-8.03#
9	A	Endosulfan I	50.000	0.000	100.0#	0	-8.44#
10	B	gamma-Chlordane	50.000	0.000	100.0#	0	-8.30#
11	B	alpha-Chlordane	50.000	0.000	100.0#	0	-8.38#
12	B	4,4'-DDE	50.000	0.000	100.0#	0	-8.57#
13	MA	Dieldrin	50.000	0.000	100.0#	0	-8.73#
14	MA	Endrin	50.000	0.000	100.0#	0	-8.97#
15	B	Endosulfan II	50.000	0.000	100.0#	0	-9.20#
16	A	4,4'-DDD	50.000	0.000	100.0#	0	-9.11#
17	MA	4,4'-DDT	50.000	0.000	100.0#	0	-9.43#
18	B	Endrin aldehyde	50.000	0.000	100.0#	0	-9.34#
19	B	Endosulfan Sulfate	50.000	0.000	100.0#	0	-9.57#
20	A	Methoxychlor	50.000	0.000	100.0#	0	-9.91#
21	B	Endrin ketone	50.000	0.000	100.0#	0	-10.07#
22		Mirex	50.000	0.000	100.0#	0	-10.54#

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