

Method Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\
 Method File : PD041423CLP.M
 Title : GC Extractables
 Last Update : Fri Apr 14 02:01:12 2023
 Response Via : Initial Calibration

Calibration Files

3 =PD074708.D 1 =PD074706.D 2 =PD074707.D
 4 =PD074709.D 5 =PD074710.D

	Compound	3	1	2	4	5	Avg	%RSD
1) SA	Tetrachloro-m-xylene	2.970	3.035	3.033	2.931	2.797	2.953 E6	3.31
2) A	alpha-BHC						0.000	-1.00
3) MA	gamma-BHC (Lindane)						0.000	-1.00
4) MA	Heptachlor						0.000	-1.00
5) MB	Aldrin						0.000	-1.00
6) B	beta-BHC						0.000	-1.00
7) B	delta-BHC						0.000	-1.00
8) B	Heptachlor epoxide						0.000	-1.00
9) A	Endosulfan I						0.000	-1.00
10) B	trans-Chlordane						0.000	-1.00
11) B	cis-Chlordane						0.000	-1.00
12) B	4,4'-DDE						0.000	-1.00
13) MA	Dieldrin						0.000	-1.00
14) MA	Endrin						0.000	-1.00
15) B	Endosulfan II						0.000	-1.00
16) A	4,4'-DDD						0.000	-1.00
17) MA	4,4'-DDT						0.000	-1.00
18) B	Endrin aldehyde						0.000	-1.00
19) B	Endosulfan Sulfate						0.000	-1.00
20) A	Methoxychlor						0.000	-1.00
21) B	Endrin ketone						0.000	-1.00
22)	Toxaphene-1	2.762	2.897	2.787	2.764	2.688	2.780 E4	2.72
23)	Toxaphene-2	3.451	3.500	3.435	3.544	3.348	3.455 E4	2.13
24)	Toxaphene-3	2.336	2.199	2.300	2.442	2.308	2.317 E4	3.75
25)	Toxaphene-4	7.702	7.406	7.533	8.068	7.949	7.732 E4	3.58
26)	Toxaphene-5	5.816	5.490	5.526	6.067	5.902	5.760 E4	4.30
27) SA	Decachlorobiphenyl	3.623	4.056	3.829	3.466	3.255	3.646 E6	8.53

Signal #2 Calibration Files

3 =PD074708.D 1 =PD074706.D 2 =PD074707.D
 4 =PD074709.D 5 =PD074710.D

	Compound	3	1	2	4	5	Avg	%RSD
1) SA	Tetrachloro-m-xylene	1.651	1.674	1.685	1.604	1.510	1.625 E6	4.39
2) A	alpha-BHC						0.000	-1.00
3) MA	gamma-BHC (Lindane)						0.000	-1.00
4) MA	Heptachlor						0.000	-1.00
5) MB	Aldrin						0.000	-1.00
6) B	beta-BHC						0.000	-1.00
7) B	delta-BHC						0.000	-1.00
8) B	Heptachlor epoxide						0.000	-1.00
9) A	Endosulfan I						0.000	-1.00
10) B	trans-Chlordane						0.000	-1.00
11) B	cis-Chlordane						0.000	-1.00
12) B	4,4'-DDE						0.000	-1.00
13) MA	Dieldrin						0.000	-1.00
14) MA	Endrin						0.000	-1.00
15) B	Endosulfan II						0.000	-1.00
16) A	4,4'-DDD						0.000	-1.00
17) MA	4,4'-DDT						0.000	-1.00
18) B	Endrin aldehyde						0.000	-1.00
19) B	Endosulfan Sulfate						0.000	-1.00
20) A	Methoxychlor						0.000	-1.00
21) B	Endrin ketone						0.000	-1.00
22)	Toxaphene-1	0.984	1.042	1.024	0.969	0.905	0.985 E4	5.43
23)	Toxaphene-2	1.085	1.119	1.093	1.121	1.062	1.096 E4	2.23

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	Compound	3	1	2	4	5	Avg		%RSD
24)	Toxaphene-3	1.215	1.299	1.188	1.224	1.133	1.212	E4	4.97
25)	Toxaphene-4	3.841	3.584	3.667	3.968	3.919	3.796	E4	4.34
26)	Toxaphene-5	5.096	4.876	5.022	5.065	4.955	5.002	E4	1.77
27) SA	Decachlorobiphenyl	1.689	1.814	1.748	1.583	1.484	1.663	E6	7.90

(#) = Out of Range