

Method Path : Z:\pestpcbsrv\HPCHEM1\ECD\_L\methods\  
 Method File : PL010422.M  
 Title : GC Extractables  
 Last Update : Wed Jan 05 01:12:56 2022  
 Response Via : Initial Calibration

## Calibration Files

50 =PL071794.D 100 =PL071792.D 75 =PL071793.D  
 25 =PL071795.D 5 =PL071796.D

|        | Compound             | 50    | 100   | 75    | 25    | 5     | Avg      | %RSD  |
|--------|----------------------|-------|-------|-------|-------|-------|----------|-------|
| 1) SA  | Tetrachloro-m-xylene | 4.901 | 4.894 | 4.880 | 4.850 | 5.618 | 5.029 E5 | 6.57  |
| 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  | gamma-Chlordane      |       |       |       |       |       | 0.000    | -1.00 |
| 11) B  | alpha-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)    | Mirex                |       |       |       |       |       | 0.000    | -1.00 |
| 23)    | Chlordane-1          | 1.807 | 1.809 | 1.821 | 1.739 | 1.757 | 1.787 E4 | 2.04  |
| 24)    | Chlordane-2          | 2.436 | 2.324 | 2.392 | 2.445 | 2.422 | 2.404 E4 | 2.04  |
| 25)    | Chlordane-3          | 6.701 | 6.710 | 6.735 | 6.392 | 6.265 | 6.561 E4 | 3.31  |
| 26)    | Chlordane-4          | 7.167 | 7.229 | 7.211 | 6.927 | 6.768 | 7.060 E4 | 2.88  |
| 27)    | Chlordane-5          | 2.434 | 2.397 | 2.431 | 2.365 | 2.345 | 2.394 E4 | 1.64  |
| 28) SA | Decachlorobiphenyl   | 7.311 | 6.725 | 6.962 | 7.657 | 8.262 | 7.383 E5 | 8.19  |

## Signal #2 Calibration Files

50 =PL071794.D 100 =PL071792.D 75 =PL071793.D  
 25 =PL071795.D 5 =PL071796.D

|        | Compound             | 50    | 100   | 75    | 25    | 5     | Avg      | %RSD  |
|--------|----------------------|-------|-------|-------|-------|-------|----------|-------|
| 1) SA  | Tetrachloro-m-xylene | 1.274 | 1.257 | 1.269 | 1.253 | 1.213 | 1.253 E6 | 1.93  |
| 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  | gamma-Chlordane      |       |       |       |       |       | 0.000    | -1.00 |
| 11) B  | alpha-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)    | Mirex                |       |       |       |       |       | 0.000    | -1.00 |

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|        | Compound           | 50    | 100   | 75    | 25    | 5     | Avg   |    | %RSD |
|--------|--------------------|-------|-------|-------|-------|-------|-------|----|------|
| -----  |                    |       |       |       |       |       |       |    |      |
| 23)    | Chlordane-1        | 5.128 | 5.138 | 5.099 | 4.935 | 4.466 | 4.953 | E4 | 5.74 |
| 24)    | Chlordane-2        | 5.420 | 5.229 | 5.369 | 5.503 | 5.543 | 5.413 | E4 | 2.28 |
| 25)    | Chlordane-3        | 2.083 | 2.101 | 2.116 | 2.007 | 1.898 | 2.041 | E5 | 4.42 |
| 26)    | Chlordane-4        | 2.488 | 2.474 | 2.498 | 2.418 | 2.292 | 2.434 | E5 | 3.50 |
| 27)    | Chlordane-5        | 4.105 | 4.063 | 4.067 | 3.953 | 3.830 | 4.004 | E4 | 2.80 |
| 28) SA | Decachlorobiphenyl | 1.207 | 1.162 | 1.184 | 1.224 | 1.253 | 1.206 | E6 | 2.93 |
| -----  |                    |       |       |       |       |       |       |    |      |

(#) = Out of Range