

**Instrument :**  
BNA\_G  
**ClientSampleId :**  
SSTD080305

mohammad  
7/9/2021 2:47:03 PM

Abundance

TIC: BG049227.D

Time-->

1,4-Dioxane-d8,S  
Pyridine-d5,S  
Bis(2-chloroethyl)amine-d8,S  
Bis(2-chloroethyl)amine,S  
1,4-Dichlorobenzene-d4,I  
2,2'-oxybis[propane]  
2,2'-oxybis[propane]  
N,N'-bis(2-chloroethyl)amine,P  
N,N'-bis(2-chloroethyl)amine,S  
Isobutanol  
2-Methyl-2-phenol  
Bis(2-chloroethyl)amine,S  
2,4-Dichlorobenzene-d4,S  
Naphthalene-d8,S  
Hexachlorobutadiene,C  
Caprolactam  
4-Chloro-3-methylphenol,C  
2-Methylphenol  
Hexachlorobutadiene,C  
2,4,5-Trichlorobenzene  
2-Chlorobenzene  
2-Nitroaniline  
2,6-Dimethylphenol-d6,S  
2,6-Dimethylphenol,S  
Acetophenone  
3-Nitrophenol  
2,4-Dichlorobenzene  
2,3,4-Trichlorobenzene  
4-Nitrophenol  
4-Nitro-2-dimethylphenylamine  
4-Nitrophenyl-phenylether  
4-Bromophenyl-phenylether  
4-Bromophenyl-phenol,C  
Pentachlorophenol,C  
Phenanthrene-d10,I  
Phenanthrene,S  
Carbazole  
Di-n-butylphthalate  
Fluoranthene  
Pyrene-d10,S  
Butylbenzylphthalate  
Chrysene-d12,I  
Chrysene,S  
Benzo(a)anthracene  
Di-n-octyl phthalate  
Benzo(a)anthracene  
Benzo(a)pyrene-d12,I  
Benzo(a)pyrene,S  
Indeno(1,2,3-cd)pyrene  
Benzo(g,h,i)perylene

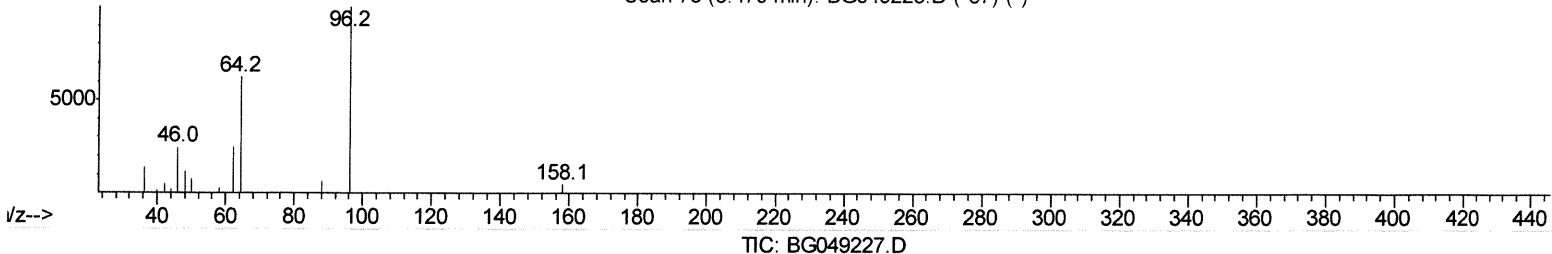
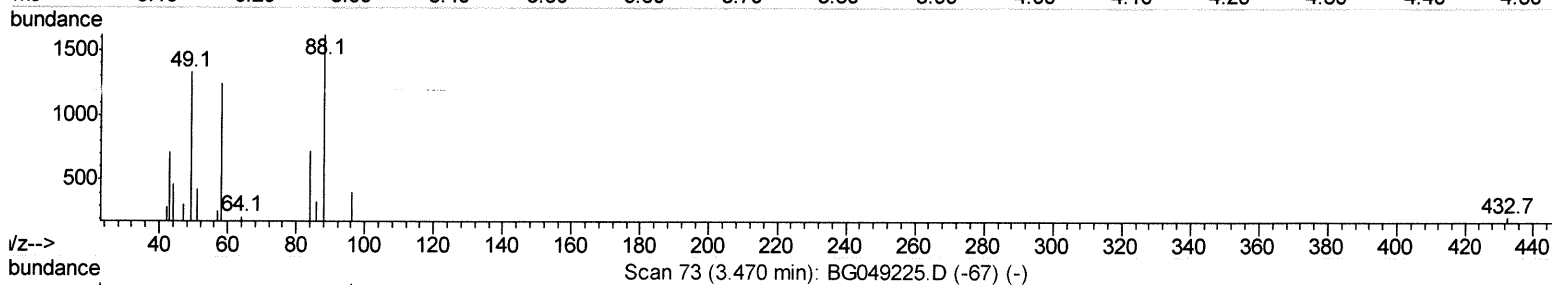
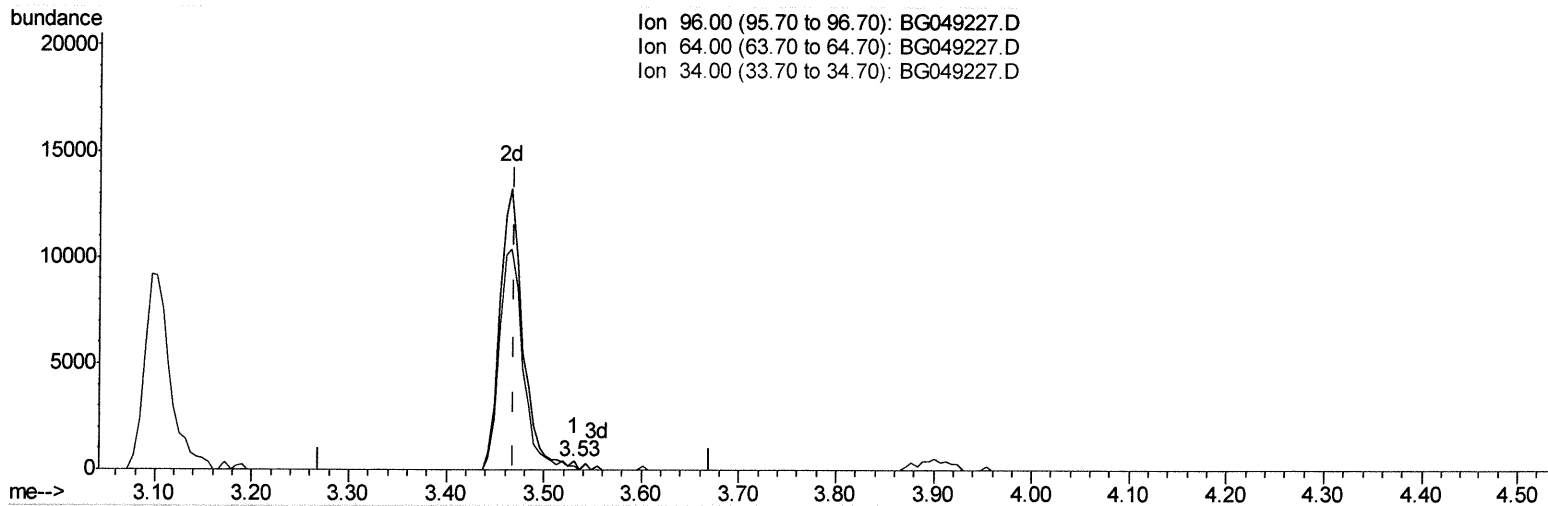
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 Data File : BG049227.D  
 Acq On : 8 Jul 2021 22:28  
 Operator : CG/JU  
 Sample : SSTD08001  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
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Manual Integrations  
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Quant Time: Jul 09 02:36:46 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG070921.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jul 09 02:34:40 2021  
 Response via : Initial Calibration



(3) 1,4-Dioxane-d8 (S)

3.530min (+0.061) 0.21ng/uL

response 145

| Ion   | Exp%  | Act%  |
|-------|-------|-------|
| 96.00 | 100   | 100   |
| 64.00 | 62.20 | 50.98 |
| 34.00 | 0.00  | 0.00  |
| 0.00  | 0.00  | 0.00  |

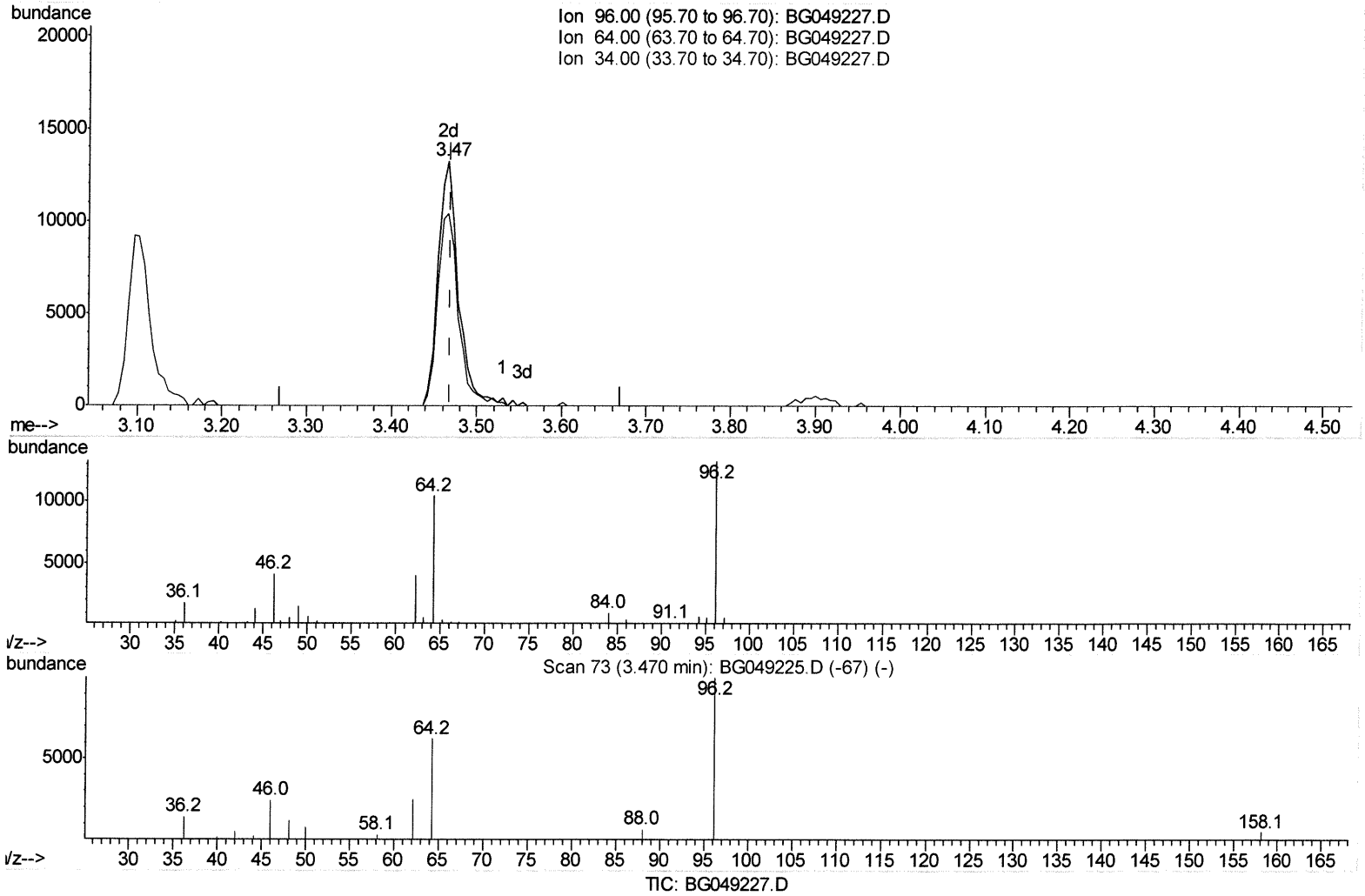
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(3) 1,4-Dioxane-d8 (S)

3.466min (-0.004) 30.90ng/uL m 07/02/21JU

response 21840

| Ion   | Exp%  | Act%   |
|-------|-------|--------|
| 96.00 | 100   | 100    |
| 64.00 | 62.20 | 78.84# |
| 34.00 | 0.00  | 0.00   |
| 0.00  | 0.00  | 0.00   |

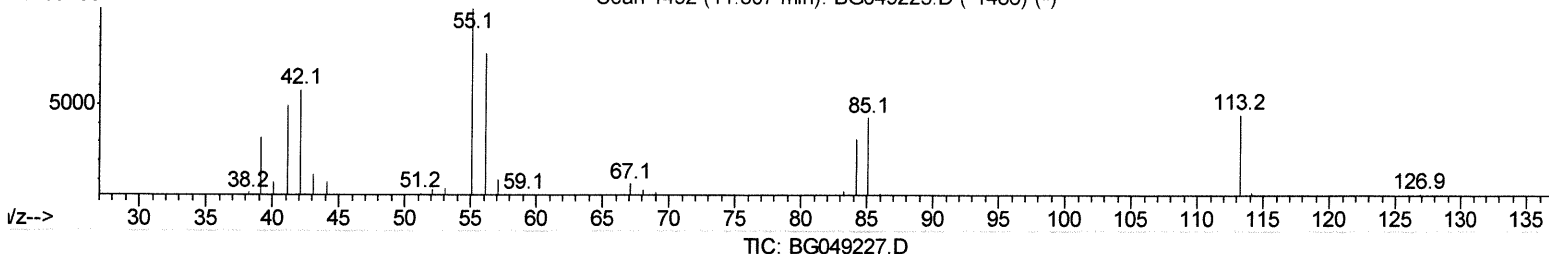
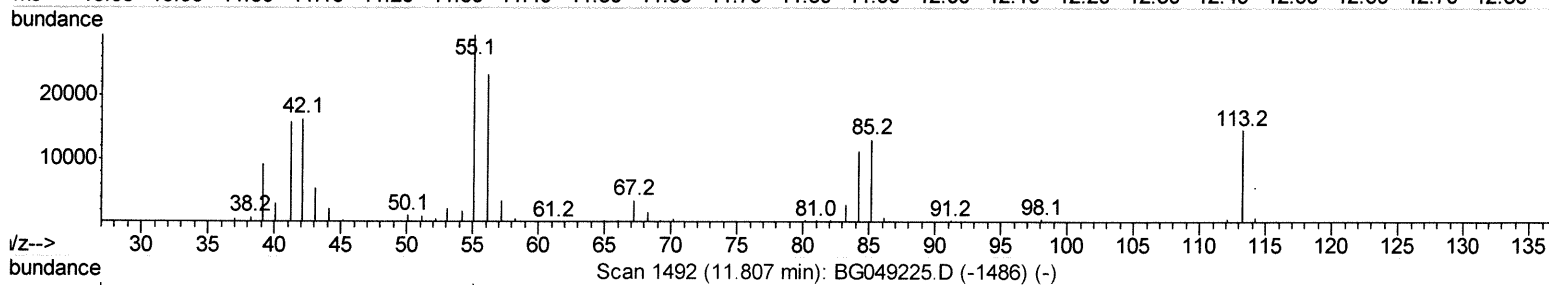
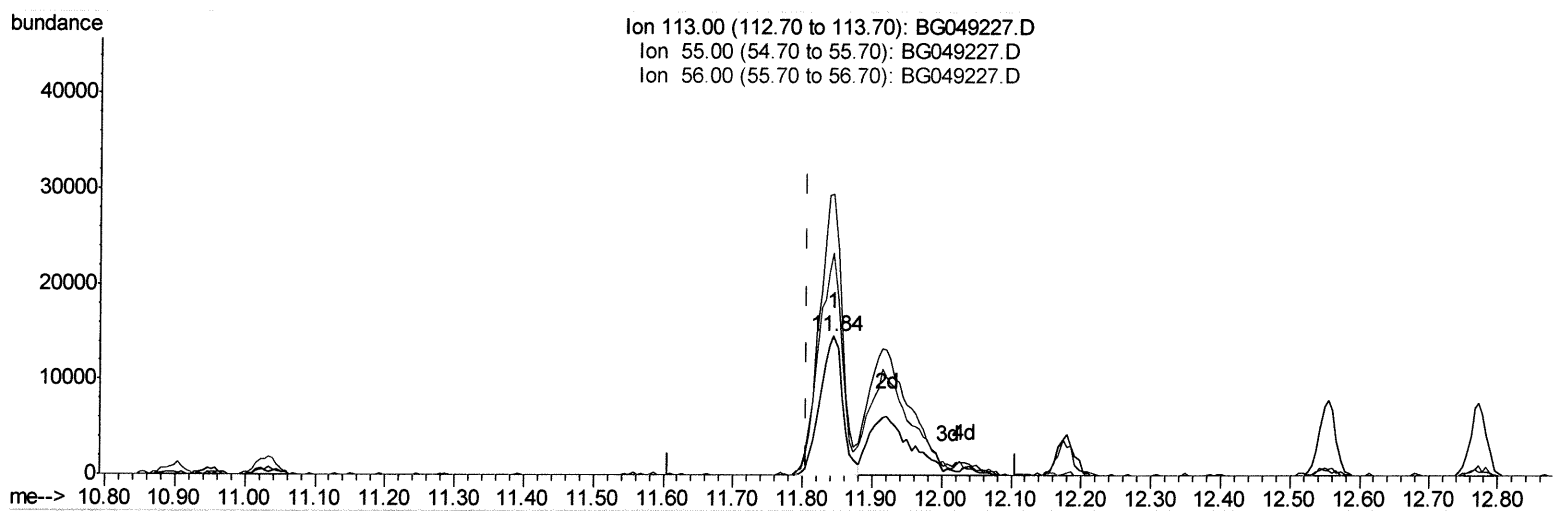
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(34) Caprolactam

11.844min (+0.037) 37.15ng/ul

response 32538

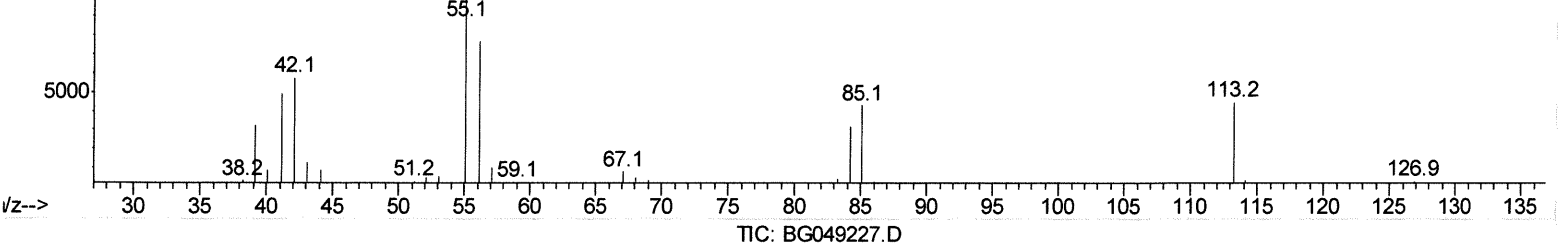
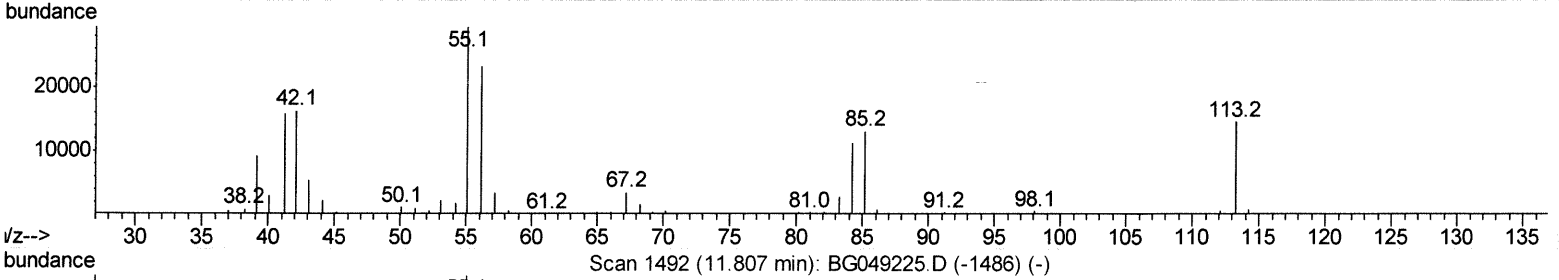
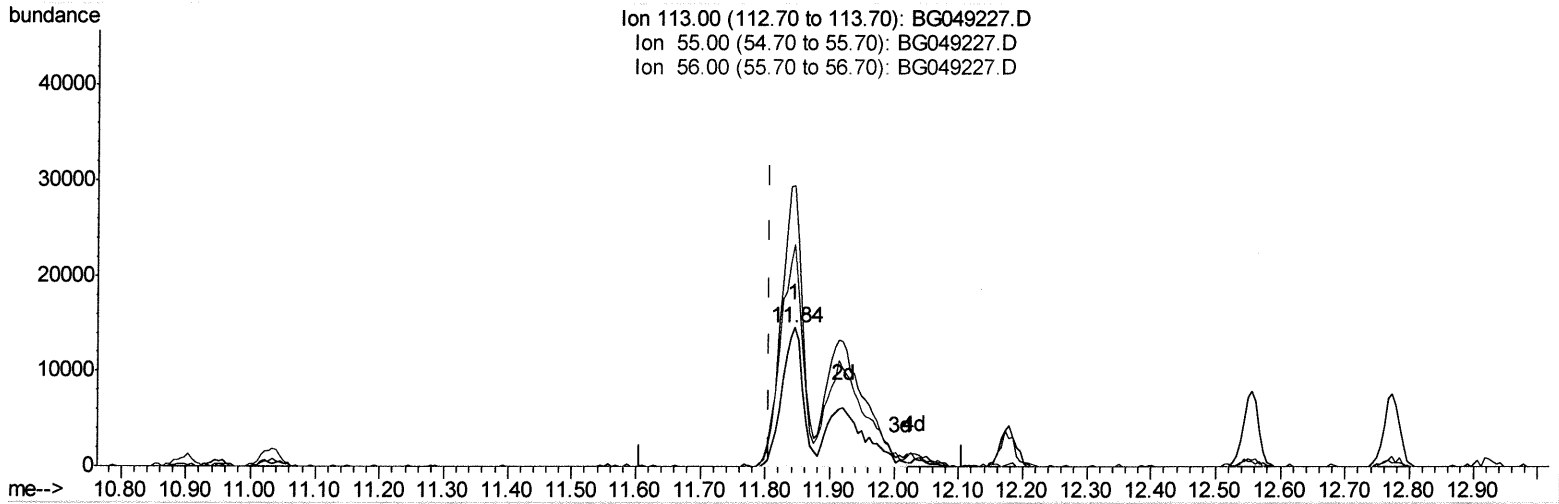
| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 223.70 | 202.32 |
| 56.00  | 171.90 | 160.03 |
| 0.00   | 0.00   | 0.00   |

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Response via : Initial Calibration



(34) Caprolactam

11.844min (+0.037) 67.05ng/ul m 07/12/21 JU

response 58732

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 223.70 | 202.32 |
| 56.00  | 171.90 | 160.03 |
| 0.00   | 0.00   | 0.00   |

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| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.08  | 152  | 30621    | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.90 | 136  | 132128   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.72 | 164  | 95125    | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.47 | 188  | 222190   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.76 | 240  | 221813   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 25.06 | 264  | 233657   | 20.00 | ng/ul | 0.01     |

#### System Monitoring Compounds

|                                |       |     |        |        |       |      |          |
|--------------------------------|-------|-----|--------|--------|-------|------|----------|
| 3) 1,4-Dioxane-d8              | 3.47  | 96  | 21840m | 30.90  | ng/uL | 0.00 | 07/12/21 |
| 4) Pyridine-d5                 | 3.88  | 84  | 163790 | 75.70  | ng/ul | 0.00 |          |
| 7) Phenol-d5                   | 7.23  | 99  | 220044 | 82.71  | ng/ul | 0.00 |          |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.40  | 67  | 124662 | 75.88  | ng/ul | 0.00 |          |
| 11) 2-Chlorophenol-d4          | 7.61  | 132 | 167968 | 90.86  | ng/ul | 0.00 |          |
| 15) 4-Methylphenol-d8          | 8.79  | 113 | 176289 | 80.78  | ng/ul | 0.00 |          |
| 21) Nitrobenzene-d5            | 9.25  | 128 | 82953  | 90.67  | ng/ul | 0.00 |          |
| 24) 2-Nitrophenol-d4           | 9.97  | 143 | 100533 | 100.37 | ng/ul | 0.00 |          |
| 28) 2,4-Dichlorophenol-d3      | 10.52 | 165 | 190008 | 77.20  | ng/ul | 0.00 |          |
| 31) 4-Chloroaniline-d4         | 11.03 | 131 | 236287 | 76.23  | ng/ul | 0.00 |          |
| 46) Dimethylphthalate-d6       | 14.12 | 166 | 580971 | 74.24  | ng/ul | 0.00 |          |
| 49) Acenaphthylene-d8          | 14.42 | 160 | 682455 | 79.42  | ng/ul | 0.00 |          |
| 54) 4-Nitrophenol-d4           | 14.93 | 143 | 97877  | 83.99  | ng/ul | 0.01 |          |
| 60) Fluorene-d10               | 15.72 | 176 | 492006 | 71.23  | ng/ul | 0.00 |          |
| 65) 4,6-Dinitro-2-methylphenol | 15.84 | 200 | 117361 | 103.36 | ng/ul | 0.01 |          |
| 73) Anthracene-d10             | 17.58 | 188 | 790479 | 77.87  | ng/ul | 0.00 |          |
| 81) Pyrene-d10                 | 19.86 | 212 | 918907 | 81.01  | ng/ul | 0.00 |          |
| 92) Benzo(a)pyrene-d12         | 24.85 | 264 | 988453 | 78.04  | ng/ul | 0.03 |          |

#### Target Compounds

|                                |       |     |        |         |        | Ovalue |
|--------------------------------|-------|-----|--------|---------|--------|--------|
| 2) 1,4-Dioxane                 | 3.50  | 88  | 23256  | 32.792  | ng/uL  | 96     |
| 5) Pyridine                    | 3.90  | 79  | 180816 | 83.664  | ng/ul  | 98     |
| 6) Benzaldehyde                | 7.21  | 77  | 133945 | 79.734  | ng/ul  | 97     |
| 8) Phenol                      | 7.26  | 94  | 218580 | 77.719  | ng/ul  | 93     |
| 10) Bis(2-Chloroethyl)ether    | 7.49  | 93  | 161460 | 78.724  | ng/ul  | 88     |
| 12) 2-Chlorophenol             | 7.64  | 128 | 164059 | 83.610  | ng/ul  | 97     |
| 13) 2-Methylphenol             | 8.52  | 108 | 168550 | 80.108  | ng/ul  | 95     |
| 14) 2,2'-oxybis(1-Chloropropan | 8.60  | 45  | 254912 | 105.198 | ng/ul  | 97     |
| 16) Acetophenone               | 8.91  | 105 | 282365 | 78.926  | ng/ul  | 96     |
| 17) N-Nitroso-di-n-propylamine | 8.90  | 70  | 147793 | 69.678  | ng/ul  | 97     |
| 18) 4-Methylphenol             | 8.85  | 108 | 173722 | 78.650  | ng/ul  | 91     |
| 19) Hexachloroethane           | 9.17  | 117 | 74636  | 81.627  | ng/ul  | 94     |
| 22) Nitrobenzene               | 9.29  | 77  | 229675 | 76.153  | ng/ul# | 95     |
| 23) Isophorone                 | 9.82  | 82  | 403469 | 64.721  | ng/ul  | 99     |
| 25) 2-Nitrophenol              | 10.01 | 139 | 97177  | 85.195  | ng/ul  | 97     |
| 26) 2,4-Dimethylphenol         | 10.06 | 107 | 213050 | 70.984  | ng/ul  | 97     |
| 27) Bis(2-Chloroethoxy)methane | 10.30 | 93  | 216471 | 68.977  | ng/ul  | 93     |
| 29) 2,4-Dichlorophenol         | 10.55 | 162 | 170061 | 73.966  | ng/ul  | 95     |
| 30) Naphthalene                | 10.95 | 128 | 528576 | 75.699  | ng/ul  | 99     |
| 32) 4-Chloroaniline            | 11.06 | 127 | 231193 | 75.845  | ng/ul  | 96     |
| 33) Hexachlorobutadiene        | 11.23 | 225 | 139311 | 63.699  | ng/ul  | 95     |
| 34) Caprolactam                | 11.84 | 113 | 58732m | 67.049  | ng/ul  | 97     |

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| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 35) 4-Chloro-3-methylphenol    | 12.18 | 107  | 192715   | 71.469 | ng/ul  | 99       |
| 36) 2-Methylnaphthalene        | 12.56 | 142  | 405539   | 75.295 | ng/ul  | 92       |
| 37) 1-Methylnaphthalene        | 12.77 | 142  | 398832   | 74.764 | ng/ul# | 95       |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.92 | 216  | 241543   | 66.704 | ng/ul# | 97       |
| 40) Hexachlorocyclopentadiene  | 12.90 | 237  | 171319   | 65.154 | ng/ul  | 98       |
| 41) 2,4,6-Trichlorophenol      | 13.16 | 196  | 161304   | 72.387 | ng/ul  | 96       |
| 42) 2,4,5-Trichlorophenol      | 13.23 | 196  | 168284   | 70.355 | ng/ul  | 88       |
| 43) 1,1'-Biphenyl              | 13.56 | 154  | 505670   | 71.840 | ng/ul  | 98       |
| 44) 2-Chloronaphthalene        | 13.60 | 162  | 397483   | 72.091 | ng/ul  | 98       |
| 45) 2-Nitroaniline             | 13.80 | 65   | 152244   | 89.386 | ng/ul  | 95       |
| 47) Dimethylphthalate          | 14.18 | 163  | 530926   | 67.170 | ng/ul# | 98       |
| 48) 2,6-Dinitrotoluene         | 14.30 | 165  | 121142   | 89.488 | ng/ul  | 87       |
| 50) Acenaphthylene             | 14.45 | 152  | 602097   | 73.068 | ng/ul  | 98       |
| 51) 3-Nitroaniline             | 14.63 | 138  | 107828   | 82.848 | ng/ul  | 98       |
| 52) Acenaphthene               | 14.79 | 153  | 440527   | 72.085 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol          | 14.83 | 184  | 80805    | 89.724 | ng/ul  | 93       |
| 55) 4-Nitrophenol              | 14.94 | 109  | 124744   | 80.820 | ng/ul  | 96       |
| 56) Dibenzofuran               | 15.12 | 168  | 617555   | 71.141 | ng/ul  | 98       |
| 57) 2,4-Dinitrotoluene         | 15.09 | 165  | 174237   | 89.476 | ng/ul# | 92       |
| 58) 2,3,4,6-Tetrachlorophenol  | 15.35 | 232  | 162302   | 71.782 | ng/ul  | 92       |
| 59) Diethylphthalate           | 15.54 | 149  | 575697   | 72.877 | ng/ul  | 98       |
| 61) Fluorene                   | 15.77 | 166  | 520442   | 70.433 | ng/ul# | 90       |
| 62) 4-Chlorophenyl-phenylether | 15.76 | 204  | 288551   | 64.376 | ng/ul  | 98       |
| 63) 4-Nitroaniline             | 15.80 | 138  | 105755   | 78.039 | ng/ul  | 86       |
| 66) 4,6-Dinitro-2-methylphenol | 15.85 | 198  | 111683   | 93.680 | ng/ul# | 99       |
| 67) N-Nitrosodiphenylamine     | 15.97 | 169  | 457241   | 74.296 | ng/ul  | 100      |
| 68) 4-Bromophenyl-phenylether  | 16.66 | 248  | 208143   | 73.656 | ng/ul  | 95       |
| 69) Hexachlorobenzene          | 16.78 | 284  | 228909   | 77.076 | ng/ul  | 95       |
| 70) Atrazine                   | 16.93 | 200  | 197983   | 71.483 | ng/ul  | 96       |
| 71) Pentachlorophenol          | 17.12 | 266  | 143973   | 76.413 | ng/ul  | 100      |
| 72) Phenanthrene               | 17.52 | 178  | 851108   | 75.441 | ng/ul  | 96       |
| 74) Anthracene                 | 17.61 | 178  | 846805   | 72.962 | ng/ul  | 97       |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.52 | 216  | 250523   | 71.917 | ng/uL  | 98       |
| 76) Pentachlorobenzene         | 15.05 | 250  | 262983   | 72.070 | ng/uL  | 99       |
| 77) Carbazole                  | 17.88 | 167  | 776266   | 78.945 | ng/ul  | 99       |
| 78) Di-n-butylphthalate        | 18.43 | 149  | 959233   | 77.943 | ng/ul  | 99       |
| 80) Fluoranthene               | 19.52 | 202  | 1070662  | 76.064 | ng/ul# | 97       |
| 82) Pyrene                     | 19.89 | 202  | 1025582  | 72.962 | ng/ul  | 98       |
| 83) Butylbenzylphthalate       | 20.76 | 149  | 426665   | 84.876 | ng/ul  | 96       |
| 84) 3,3'-Dichlorobenzidine     | 21.65 | 252  | 409752   | 75.417 | ng/ul  | 100      |
| 85) Benzo(a)anthracene         | 21.74 | 228  | 1062728  | 72.514 | ng/ul  | 97       |
| 86) Bis(2-ethylhexyl)phthalate | 21.64 | 149  | 624163   | 85.174 | ng/ul  | 99       |
| 87) Chrysene                   | 21.81 | 228  | 1016307  | 72.521 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate       | 22.88 | 149  | 1050673  | 82.610 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene       | 24.02 | 252  | 1162302  | 75.209 | ng/ul  | 98       |
| 91) Benzo(k)fluoranthene       | 24.09 | 252  | 1097417  | 72.259 | ng/ul  | 95       |
| 93) Benzo(a)pyrene             | 24.92 | 252  | 1017315  | 74.932 | ng/ul  | 97       |
| 94) Indeno(1,2,3-cd)pyrene     | 28.85 | 276  | 1337423  | 76.581 | ng/ul  | 96       |
| 95) Dibenzo(a,h)anthracene     | 28.92 | 278  | 1091783  | 76.894 | ng/ul  | 96       |
| 96) Benzo(g,h,i)perylene       | 30.05 | 276  | 1098452  | 77.599 | ng/ul  | 94       |

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|       |  |  |  |  |  |  |
|-------|--|--|--|--|--|--|
| ----- |  |  |  |  |  |  |
| ----- |  |  |  |  |  |  |

(#) = qualifier out of range (m) = manual integration (+) = signals summed