

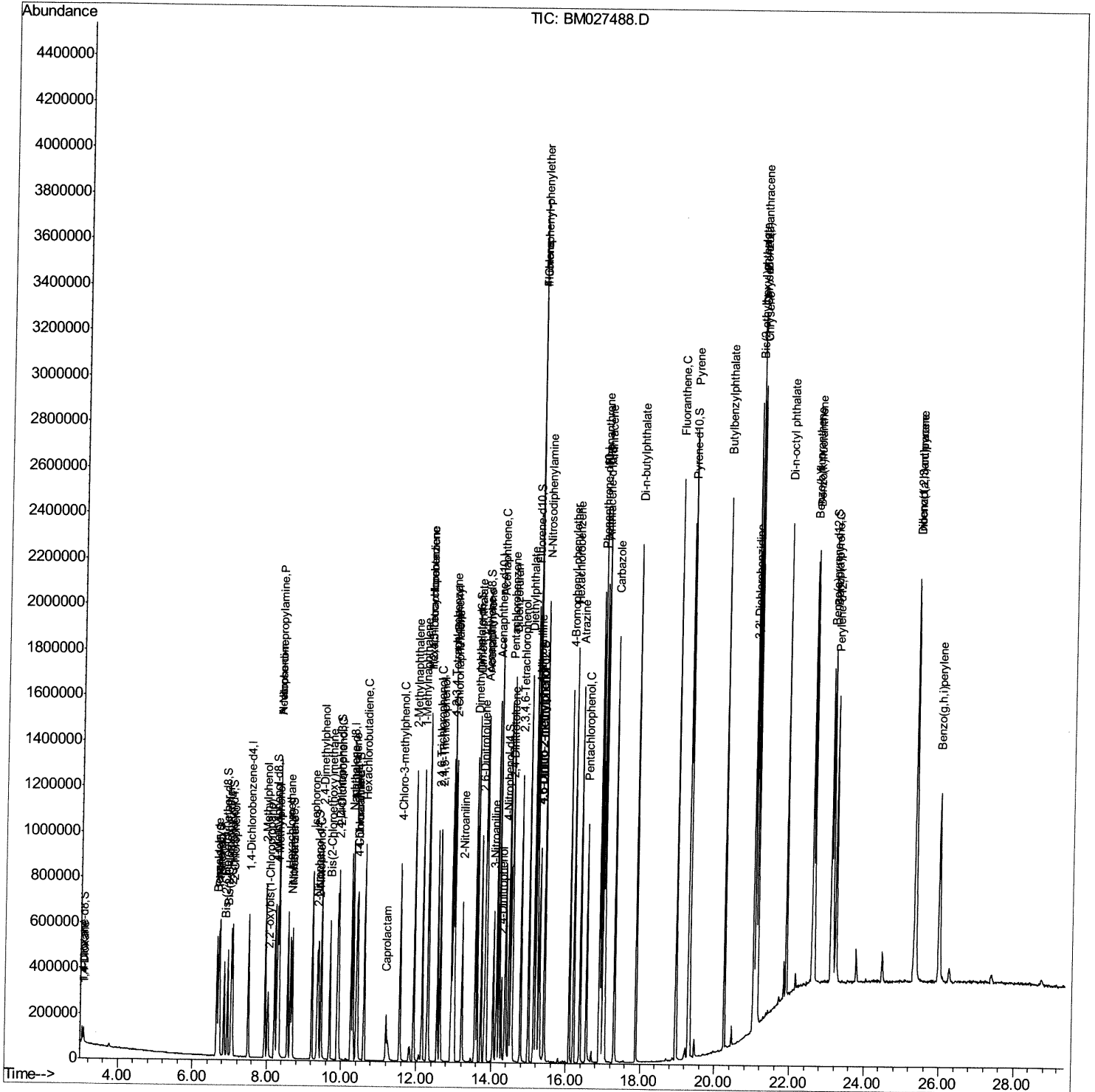
```
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091920\  
Data File : BM027488.D  
Acq On    : 19 Sep 2020 07:41  
Operator  : JU/CG  
Sample    : SSTDCCC020EC  
Misc      :  
ALS Vial  : 25      Sample Multiplier: 1
```

**Instrument :**  
BNA\_M  
**LabSampleId :**  
SSTD02024

## Manual Integrations

mohammad  
9/21/2020 10:48:13 AM

Quant Time: Sep 19 08:15:38 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM090420MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Sep 19 07:00:16 2020  
Response via : Initial Calibration



# Quantitation Report (Qedit)

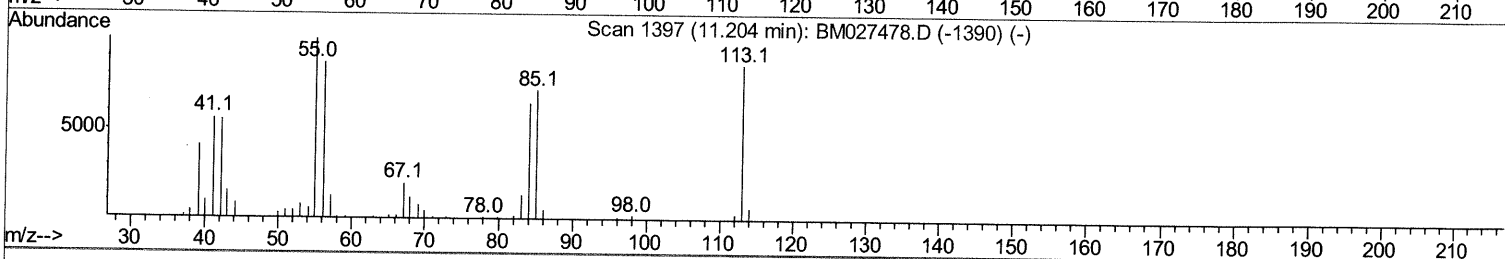
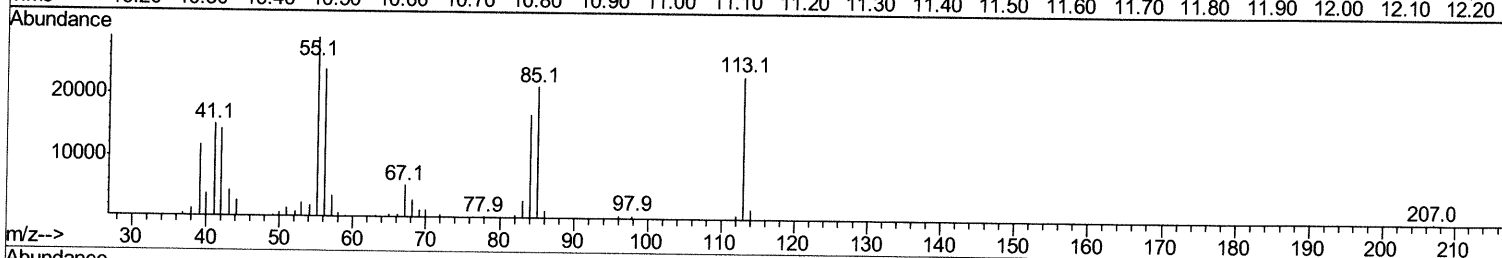
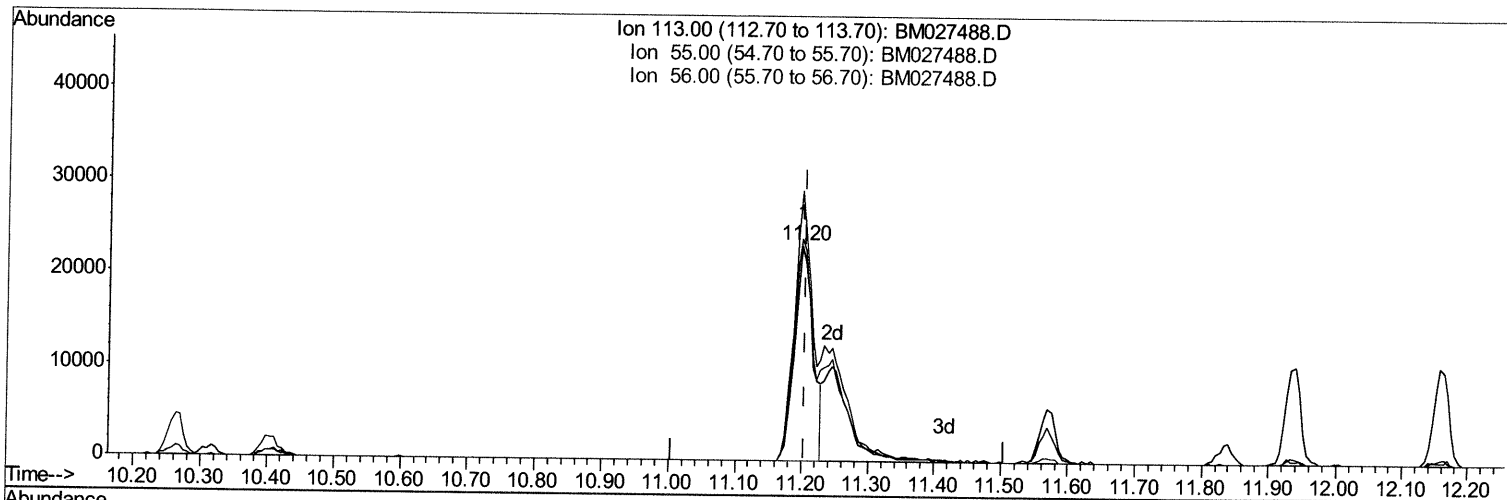
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TIC: BM027488.D

(32) Caprolactam

11.198min (-0.006) 12.17ng/ul

response 44800

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 125.60 | 125.63 |
| 56.00  | 104.60 | 103.47 |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

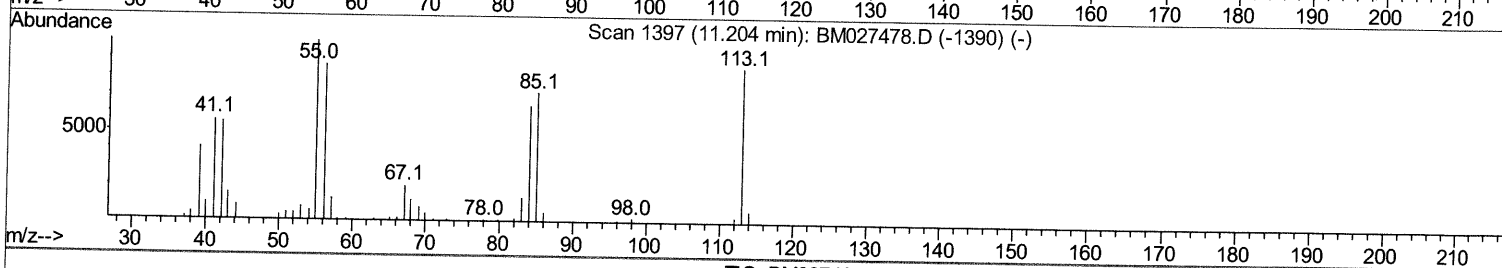
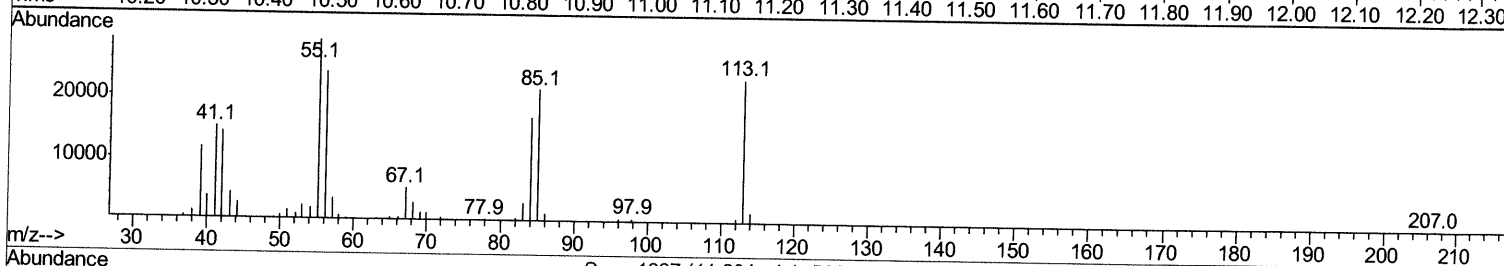
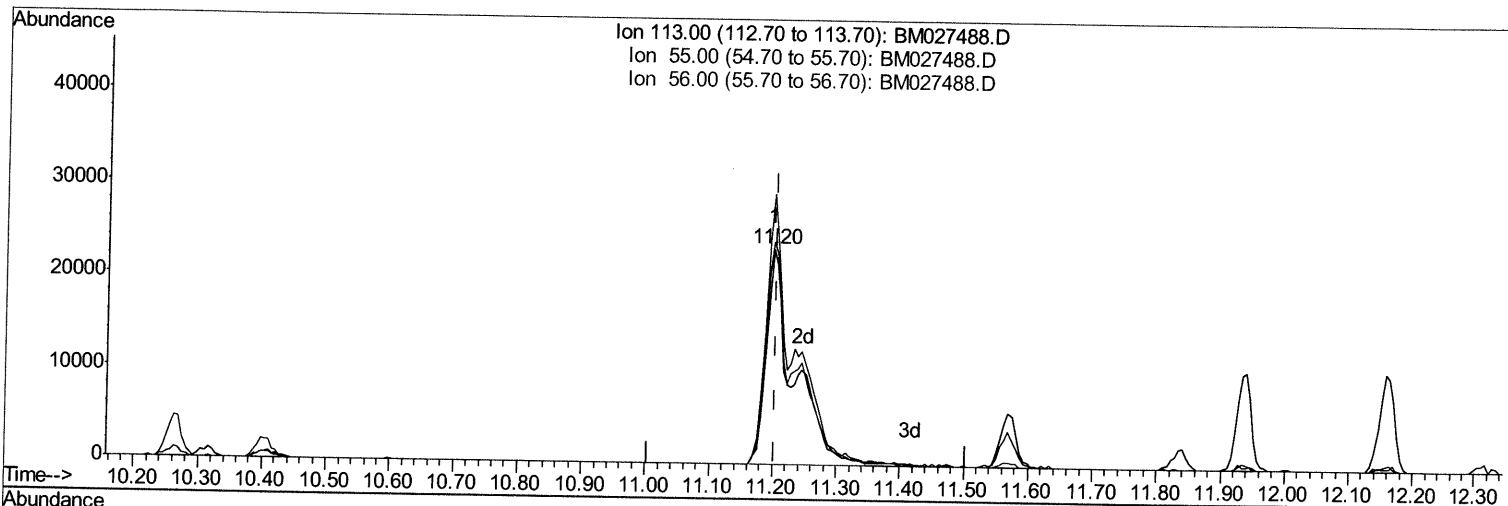
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**Manual Integrations**  
**APPROVED**

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TIC: BM027488.D

(32) Caprolactam

11.198min (-0.006) 19.35ng/ul m 09/21/20 JU

response 71234

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 125.60 | 125.63 |
| 56.00  | 104.60 | 103.47 |
| 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 | 7.49  | 152  | 160916   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.26 | 136  | 656382   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.15 | 164  | 495985   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.89 | 188  | 1153165  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.10 | 240  | 1002497  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.23 | 264  | 868425   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.05  | 96  | 34560   | 9.56  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.69  | 99  | 269736  | 20.34 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.85  | 67  | 184268  | 20.92 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.03  | 132 | 211177  | 21.74 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.22  | 113 | 220697  | 20.40 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.64  | 128 | 106321  | 21.52 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.36  | 143 | 123887  | 22.86 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.89  | 165 | 255425  | 21.90 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.40 | 131 | 283076  | 20.53 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.57 | 166 | 821944  | 20.83 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.83 | 160 | 967147  | 21.69 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.37 | 143 | 119500  | 17.19 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.14 | 176 | 734199  | 20.77 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.27 | 200 | 136665  | 18.52 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.99 | 188 | 1141732 | 20.94 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.29 | 212 | 1218648 | 27.29 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.10 | 264 | 941566  | 20.21 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |        | Ovalue |     |
|--------------------------------|-------|-----|--------|--------|--------|-----|
| 2) 1,4-Dioxane                 | 3.09  | 88  | 38446  | 8.882  | ng/uL  | 96  |
| 4) Benzaldehyde                | 6.65  | 77  | 191789 | 20.060 | ng/ul  | 95  |
| 6) Phenol                      | 6.72  | 94  | 296514 | 21.035 | ng/ul  | 92  |
| 8) Bis(2-Chloroethyl)ether     | 6.93  | 93  | 239655 | 21.168 | ng/ul  | 98  |
| 10) 2-Chlorophenol             | 7.07  | 128 | 222321 | 21.947 | ng/ul  | 97  |
| 11) 2-Methylphenol             | 7.95  | 108 | 214910 | 20.656 | ng/ul  | 98  |
| 12) 2,2'-oxybis(1-Chloropropan | 8.03  | 45  | 178841 | 20.948 | ng/ul  | 97  |
| 14) Acetophenone               | 8.32  | 105 | 383581 | 21.127 | ng/ul  | 96  |
| 15) N-Nitroso-di-n-propylamine | 8.31  | 70  | 234593 | 22.582 | ng/ul  | 99  |
| 16) 4-Methylphenol             | 8.27  | 108 | 235141 | 20.705 | ng/ul  | 95  |
| 17) Hexachloroethane           | 8.56  | 117 | 107035 | 21.237 | ng/ul  | 99  |
| 20) Nitrobenzene               | 8.68  | 77  | 330013 | 20.450 | ng/ul  | 98  |
| 21) Isophorone                 | 9.22  | 82  | 590015 | 21.542 | ng/ul  | 99  |
| 23) 2-Nitrophenol              | 9.39  | 139 | 133999 | 22.777 | ng/ul  | 98  |
| 24) 2,4-Dimethylphenol         | 9.46  | 107 | 292175 | 21.689 | ng/ul  | 99  |
| 25) Bis(2-Chloroethoxy)methane | 9.70  | 93  | 338634 | 21.763 | ng/ul  | 99  |
| 27) 2,4-Dichlorophenol         | 9.92  | 162 | 254573 | 21.924 | ng/ul  | 98  |
| 28) Naphthalene                | 10.32 | 128 | 723231 | 20.737 | ng/ul  | 100 |
| 30) 4-Chloroaniline            | 10.43 | 127 | 298707 | 21.499 | ng/ul  | 99  |
| 31) Hexachlorobutadiene        | 10.61 | 225 | 192512 | 21.020 | ng/ul  | 98  |
| 32) Caprolactam                | 11.20 | 113 | 71234m | 19.348 | ng/ul  | 99  |
| 33) 4-Chloro-3-methylphenol    | 11.57 | 107 | 270155 | 21.844 | ng/ul  | 99  |
| 34) 2-Methylnaphthalene        | 11.94 | 142 | 573659 | 21.756 | ng/ul  | 96  |

09/21/20 JU

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| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.16 | 142  | 542406   | 20.671 | ng/ul  | 99       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.32 | 216  | 368578   | 22.107 | ng/ul  | 99       |
| 38) Hexachlorocyclopentadiene  | 12.30 | 237  | 228727   | 20.966 | ng/ul  | 99       |
| 39) 2,4,6-Trichlorophenol      | 12.57 | 196  | 236706   | 23.346 | ng/ul  | 98       |
| 40) 2,4,5-Trichlorophenol      | 12.64 | 196  | 254901   | 23.277 | ng/ul  | 99       |
| 41) 1,1'-Biphenyl              | 12.97 | 154  | 804672   | 21.571 | ng/ul  | 99       |
| 42) 2-Chloronaphthalene        | 13.01 | 162  | 635714   | 20.941 | ng/ul  | 98       |
| 43) 2-Nitroaniline             | 13.22 | 65   | 198395   | 22.258 | ng/ul  | 97       |
| 45) Dimethylphthalate          | 13.62 | 163  | 862882   | 20.747 | ng/ul  | 99       |
| 46) 2,6-Dinitrotoluene         | 13.73 | 165  | 175991   | 22.192 | ng/ul  | 93       |
| 48) Acenaphthylene             | 13.86 | 152  | 967974   | 21.507 | ng/ul  | 98       |
| 49) 3-Nitroaniline             | 14.05 | 138  | 146447   | 20.731 | ng/ul  | 99       |
| 50) Acenaphthene               | 14.21 | 153  | 678175   | 20.982 | ng/ul  | 99       |
| 51) 2,4-Dinitrophenol          | 14.27 | 184  | 75534    | 15.227 | ng/ul# | 88       |
| 53) 4-Nitrophenol              | 14.38 | 109  | 126719   | 17.495 | ng/ul  | 99       |
| 54) Dibenzofuran               | 14.54 | 168  | 1015064  | 21.002 | ng/ul  | 100      |
| 55) 2,4-Dinitrotoluene         | 14.52 | 165  | 256283   | 20.965 | ng/ul  | 95       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.78 | 232  | 234203   | 21.580 | ng/ul# | 97       |
| 57) Diethylphthalate           | 14.99 | 149  | 891167   | 20.548 | ng/ul  | 100      |
| 59) Fluorene                   | 15.20 | 166  | 857006   | 20.631 | ng/ul  | 98       |
| 60) 4-Chlorophenyl-phenylether | 15.20 | 204  | 466172   | 20.506 | ng/ul  | 98       |
| 61) 4-Nitroaniline             | 15.22 | 138  | 152838   | 18.640 | ng/ul  | 98       |
| 64) 4,6-Dinitro-2-methylphenol | 15.29 | 198  | 141677   | 17.732 | ng/ul  | 93       |
| 65) N-Nitrosodiphenylamine     | 15.42 | 169  | 741815   | 22.812 | ng/ul  | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.10 | 248  | 303175   | 22.711 | ng/ul  | 97       |
| 67) Hexachlorobenzene          | 16.21 | 284  | 351606   | 21.757 | ng/ul  | 97       |
| 68) Atrazine                   | 16.38 | 200  | 287685   | 20.421 | ng/ul  | 98       |
| 69) Pentachlorophenol          | 16.55 | 266  | 171257   | 17.268 | ng/ul  | 97       |
| 70) Phenanthrene               | 16.93 | 178  | 1383950  | 20.956 | ng/ul  | 100      |
| 72) Anthracene                 | 17.03 | 178  | 1415069  | 21.033 | ng/ul  | 100      |
| 73) 1,2,3,4-Tetrachlorobenzene | 12.93 | 216  | 370496   | 24.408 | ng/uL  | 99       |
| 74) Pentachlorobenzene         | 14.47 | 250  | 382227   | 22.834 | ng/uL  | 97       |
| 75) Carbazole                  | 17.30 | 167  | 1154276  | 20.269 | ng/ul  | 100      |
| 76) Di-n-butylphthalate        | 17.89 | 149  | 1524427  | 21.554 | ng/ul  | 99       |
| 77) Fluoranthene               | 18.96 | 202  | 1567606  | 19.264 | ng/ul  | 99       |
| 80) Perylene                   | 19.33 | 202  | 1617415  | 26.670 | ng/ul  | 99       |
| 81) Butylbenzylphthalate       | 20.25 | 149  | 598275   | 25.774 | ng/ul  | 99       |
| 82) 3,3'-Dichlorobenzidine     | 21.02 | 252  | 423403   | 19.231 | ng/ul  | 97       |
| 83) Benzo(a)anthracene         | 21.08 | 228  | 1392246  | 21.434 | ng/ul  | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.04 | 149  | 862383   | 23.847 | ng/ul  | 99       |
| 85) Chrysene                   | 21.13 | 228  | 1325959  | 20.660 | ng/ul  | 99       |
| 87) Di-n-octyl phthalate       | 21.89 | 149  | 1311517  | 24.320 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 22.60 | 252  | 1233659  | 21.517 | ng/ul  | 99       |
| 89) Benzo(k)fluoranthene       | 22.64 | 252  | 1197565  | 20.834 | ng/ul  | 100      |
| 91) Benzo(a)pyrene             | 23.14 | 252  | 1032782  | 19.307 | ng/ul  | 99       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.35 | 276  | 1240515  | 17.665 | ng/ul  | 99       |
| 93) Dibenzo(a,h)anthracene     | 25.36 | 278  | 1023877  | 17.238 | ng/ul  | 99       |
| 94) Benzo(a,h,i)perylene       | 25.99 | 276  | 1017114  | 17.329 | ng/ul  | 99       |

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