

(QT Reviewed)

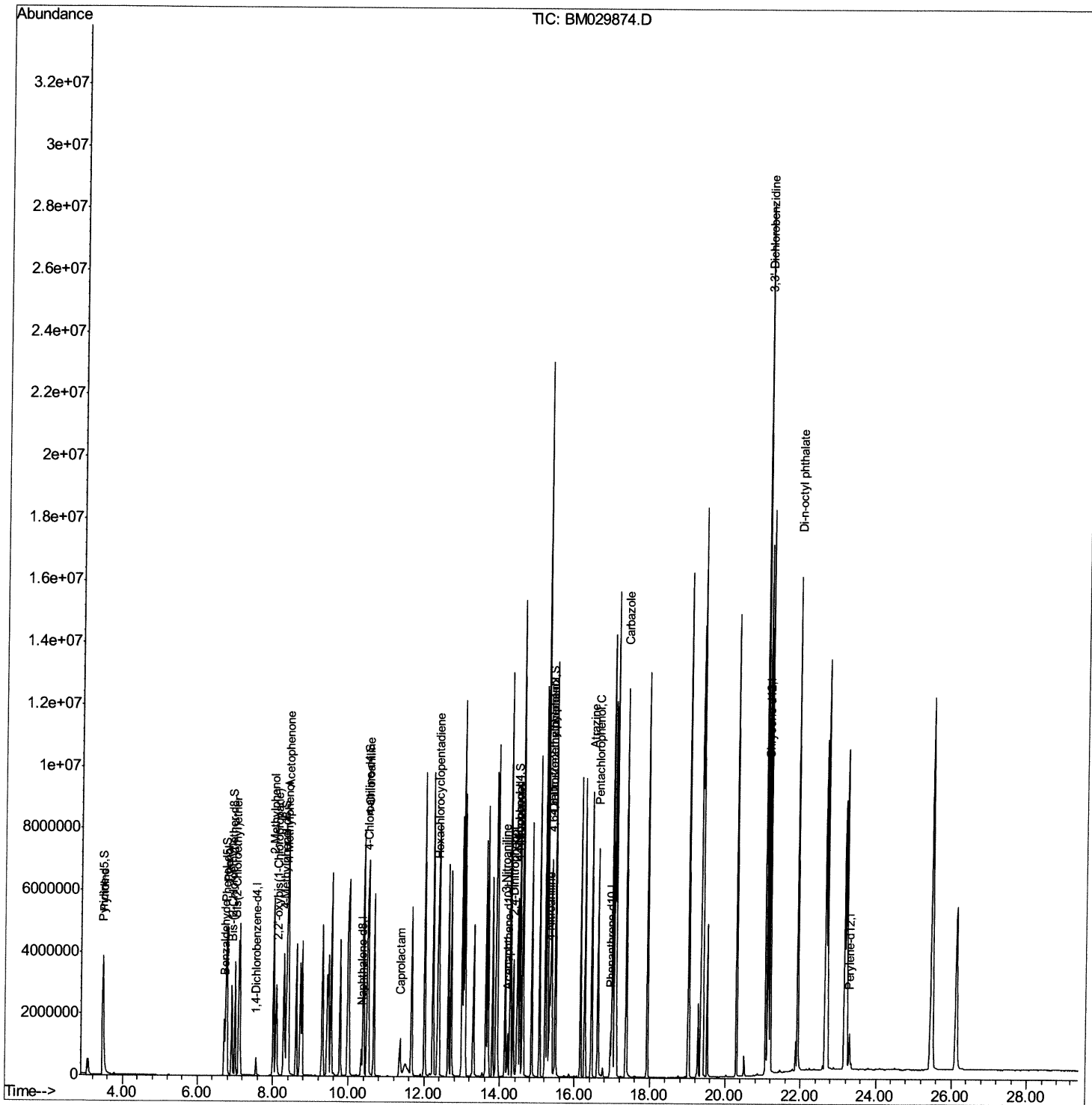
```
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050721\  
Data File : BM029874.D  
Acq On    : 07 May 2021 13:12  
Operator  : CG/JU  
Sample    : SSTD16016  
Misc      :  
ALS Vial  : 7      Sample Multiplier: 1
```

**Instrument :**  
BNA\_M  
**ClientSampleId :**  
SSTD160016

## Manual Integrations

mohammad  
5/10/2021 11:44:56 AM

Quant Time: May 07 13:48:50 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri May 07 12:05:00 2021  
Response via : Initial Calibration





## Quantitation Report (Qedit)

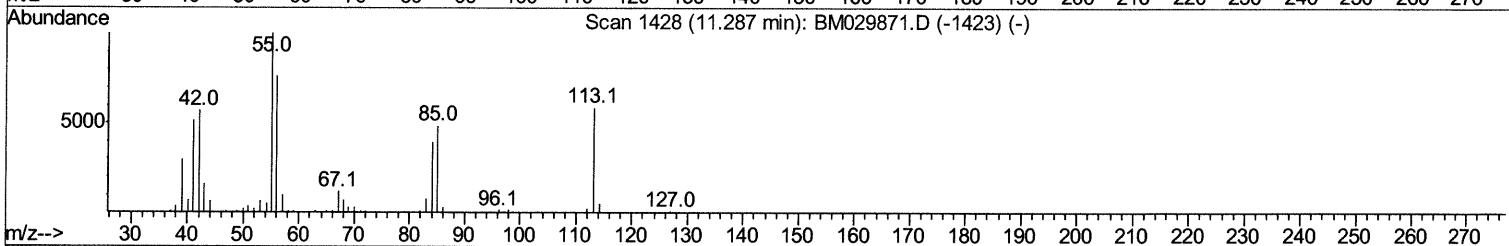
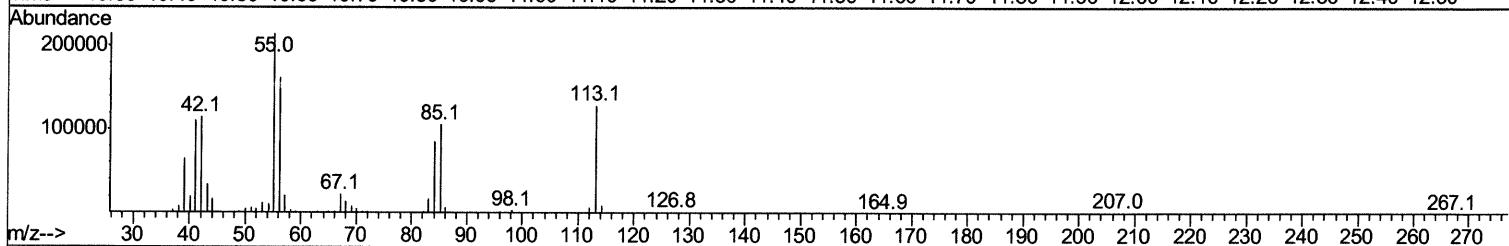
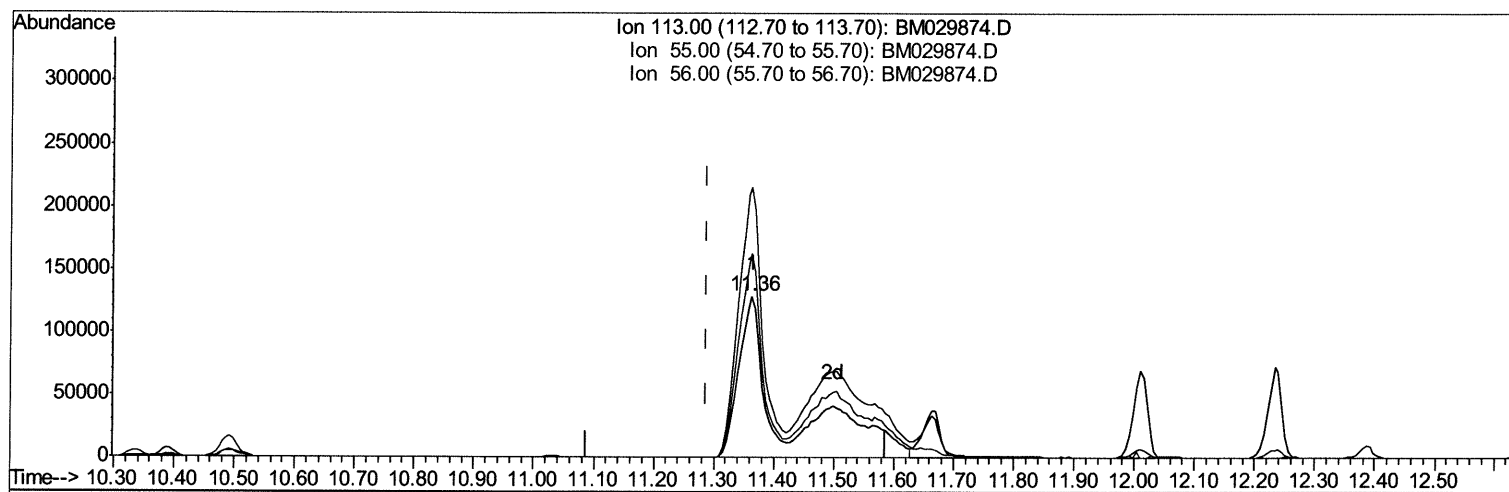
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Manual Integrations  
APPROVED

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Quant Time: May 07 13:42:35 2021  
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Quant Title : SVOA CALIBRATION  
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TIC: BM029874.D

(34) Caprolactam

11.363min (+0.076) 178.21ng/ul m 05/11/21JU

response 658225

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 173.60 | 167.57 |
| 56.00  | 131.80 | 126.43 |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

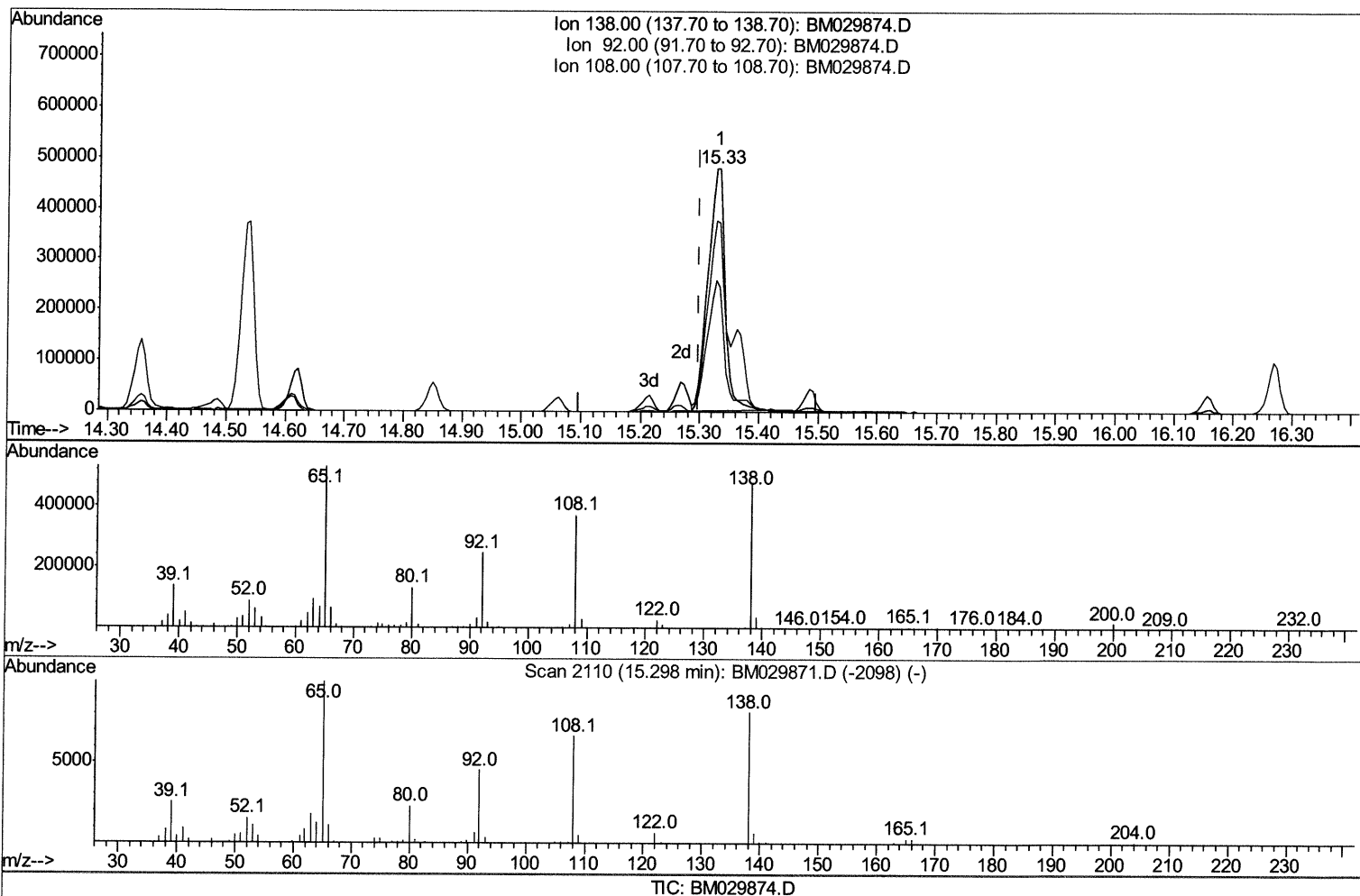
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(63) 4-Nitroaniline

15.333min (+0.035) 148.33ng/ul

response 958389

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 56.00 | 51.33 |
| 108.00 | 81.60 | 77.85 |
| 0.00   | 0.00  | 0.00  |

# Quantitation Report (Qedit)

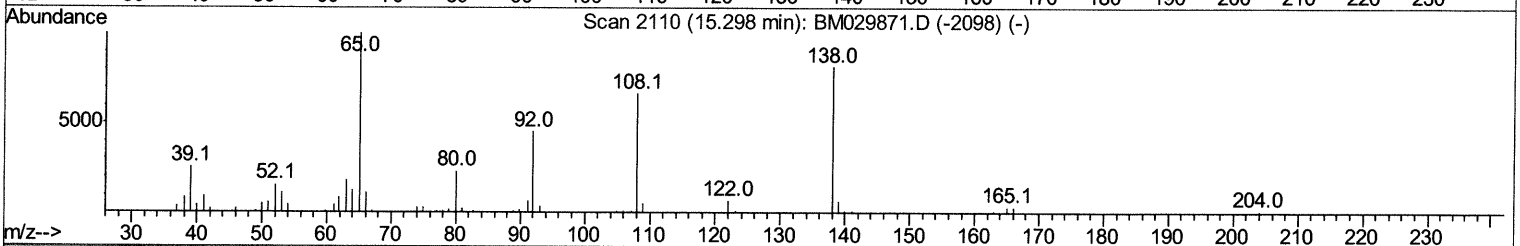
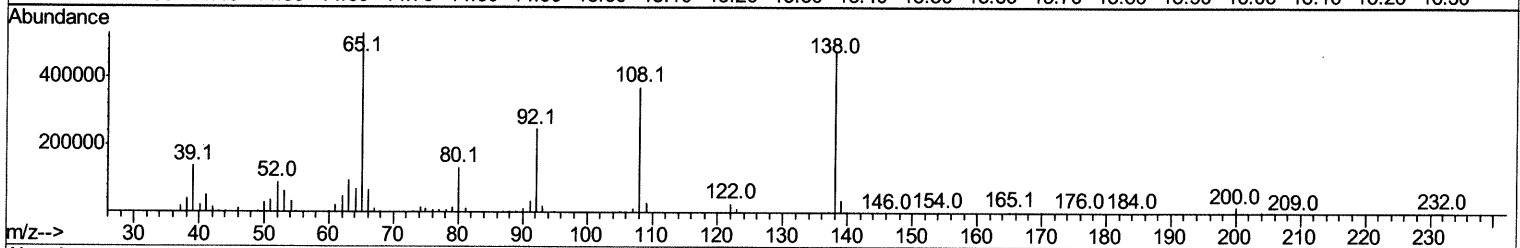
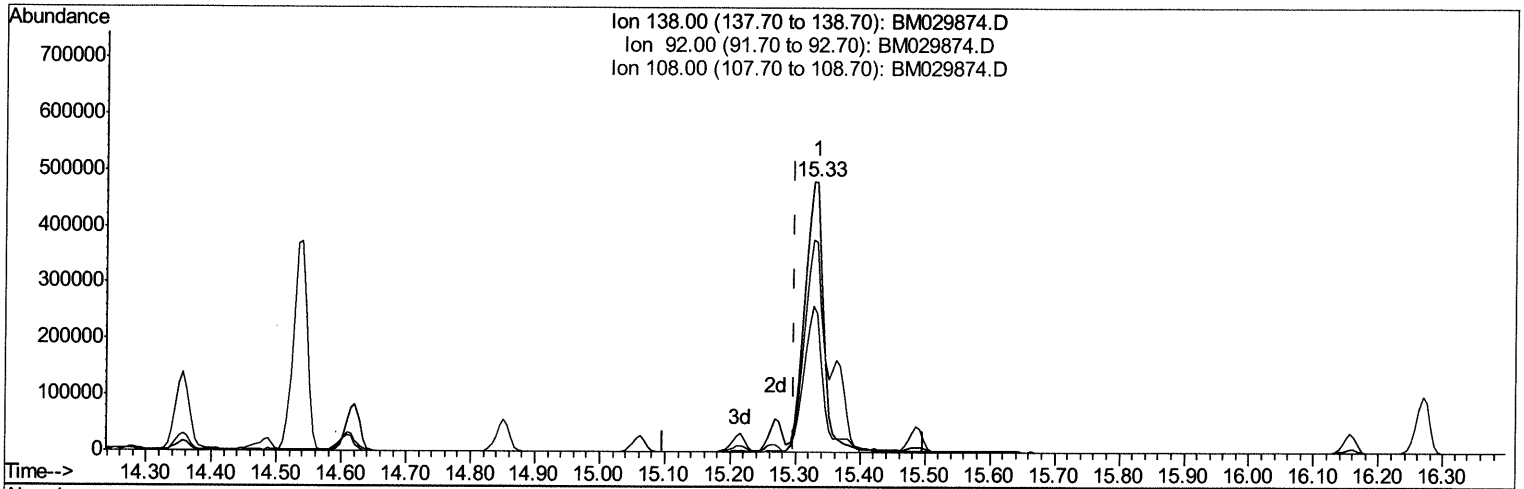
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 Operator : CG/JU  
 Sample : SSTD16016  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

**Instrument :**  
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**ClientSampleId :**  
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 Response via : Initial Calibration



TIC: BM029874.D

(63) 4-Nitroaniline

15.333min (+0.035) 165.06ng/ul m 05/10/2021

response 1066533

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 56.00 | 51.33 |
| 108.00 | 81.60 | 77.85 |
| 0.00   | 0.00  | 0.00  |

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 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.56  | 152  | 152779   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.34 | 136  | 711511   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.21 | 164  | 462973   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 16.96 | 188  | 935594   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.15 | 240  | 935744   | 20.00 | ng/ul | 0.01     |
| 88) Perylene-d12          | 23.30 | 264  | 733015   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |        |       |       |
|--------------------------------|-------|-----|---------|--------|-------|-------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96  | 0d      | 0.00   | ng/uL |       |
| 4) Pyridine-d5                 | 3.46  | 84  | 1801241 | 185.27 | ng/ul | -0.02 |
| 7) Phenol-d5                   | 6.76  | 99  | 2513138 | 178.34 | ng/ul | 0.02  |
| 9) Bis-(2-Chloroethyl)ether-d  | 6.91  | 67  | 1429880 | 169.50 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4          | 0.00  | 132 | 0d      | 0.00   | ng/ul |       |
| 15) 4-Methylphenol-d8          | 8.29  | 113 | 2027639 | 174.14 | ng/ul | 0.02  |
| 21) Nitrobenzene-d5            | 0.00  | 128 | 0d      | 0.00   | ng/ul |       |
| 24) 2-Nitrophenol-d4           | 0.00  | 143 | 0d      | 0.00   | ng/ul |       |
| 28) 2,4-Dichlorophenol-d3      | 0.00  | 165 | 0d      | 0.00   | ng/ul |       |
| 31) 4-Chloroaniline-d4         | 10.49 | 131 | 2889381 | 173.77 | ng/ul | 0.01  |
| 46) Dimethylphthalate-d6       | 0.00  | 166 | 0d      | 0.00   | ng/ul |       |
| 49) Acenaphthylene-d8          | 0.00  | 160 | 0d      | 0.00   | ng/ul |       |
| 54) 4-Nitrophenol-d4           | 14.47 | 143 | 1103427 | 205.29 | ng/ul | 0.03  |
| 60) Fluorene-d10               | 0.00  | 176 | 0d      | 0.00   | ng/ul |       |
| 65) 4,6-Dinitro-2-methylphenol | 15.36 | 200 | 1151377 | 202.27 | ng/ul | 0.02  |
| 73) Anthracene-d10             | 0.00  | 188 | 0d      | 0.00   | ng/ul |       |
| 81) Pyrene-d10                 | 0.00  | 212 | 0d      | 0.00   | ng/ul |       |
| 92) Benzo(a)pyrene-d12         | 0.00  | 264 | 0d      | 0.00   | ng/ul |       |

## Target Compounds

|                                |       |     |           |         |         | Ovalue      |
|--------------------------------|-------|-----|-----------|---------|---------|-------------|
| 5) Pyridine                    | 3.48  | 79  | 1828311   | 184.395 | ng/ul   | 92          |
| 6) Benzaldehyde                | 6.71  | 77  | 677283    | 92.518  | ng/ul   | 98          |
| 8) Phenol                      | 6.79  | 94  | 2514304   | 175.213 | ng/ul   | 99          |
| 10) Bis(2-Chloroethyl)ether    | 7.00  | 93  | 1880509   | 166.705 | ng/ul   | 97          |
| 13) 2-Methylphenol             | 8.02  | 108 | 1910433   | 174.206 | ng/ul   | 98          |
| 14) 2,2'-oxybis(1-Chloropropan | 8.10  | 45  | 2675599   | 166.324 | ng/ul   | 100         |
| 16) Acetophenone               | 8.40  | 105 | 3152491   | 170.476 | ng/ul   | 97          |
| 18) 4-Methylphenol             | 8.37  | 108 | 2091924   | 173.867 | ng/ul   | 99          |
| 32) 4-Chloroaniline            | 10.52 | 127 | 3007949   | 175.620 | ng/ul   | 99          |
| 34) Caprolactam                | 11.36 | 113 | 658225m>  | 178.207 | ng/ul>  | 05/11/21 JU |
| 40) Hexachlorocyclopentadiene  | 12.36 | 237 | 1516472   | 211.501 | ng/ul   | 98          |
| 51) 3-Nitroaniline             | 14.14 | 138 | 1208151   | 179.970 | ng/ul   | 94          |
| 53) 2,4-Dinitrophenol          | 14.36 | 184 | 816142    | 211.296 | ng/ul   | 96          |
| 55) 4-Nitrophenol              | 14.49 | 109 | 817997    | 202.506 | ng/ul   | 93          |
| 63) 4-Nitroaniline             | 15.33 | 138 | 1066533m> | 165.064 | ng/ul > | 05/11/21 JU |
| 66) 4,6-Dinitro-2-methylphenol | 15.38 | 198 | 1119275   | 196.909 | ng/ul#  | 94          |
| 70) Atrazine                   | 16.46 | 200 | 1846100   | 170.844 | ng/ul   | 99          |
| 71) Pentachlorophenol          | 16.62 | 266 | 1324862   | 187.618 | ng/ul   | 100         |
| 77) Carbazole                  | 17.37 | 167 | 8240540   | 163.664 | ng/ul   | 98          |
| 84) 3,3'-Dichlorobenzidine     | 21.07 | 252 | 2844778   | 147.490 | ng/ul   | 98          |
| 89) Di-n-octyl phthalate       | 21.93 | 149 | 10255466  | 196.917 | ng/ul   | 100         |

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Response via : Initial Calibration

| Internal Standards | R.T. | QI | on | Response | Conc | Units | Dev | (Min) |
|--------------------|------|----|----|----------|------|-------|-----|-------|
|--------------------|------|----|----|----------|------|-------|-----|-------|

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