

Data File : BN004689.D

Operator : JU/SJ

Sample : SSTDCCC020

Misc :

ALS Vial : 2 Sample Multiplier: 1

Quant Time: Mar 09 02:43:54 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN030719MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Mar 08 00:36:35 2019

Response via : Initial Calibration

**Instrument :**

BNA N

**LabSampleId :**

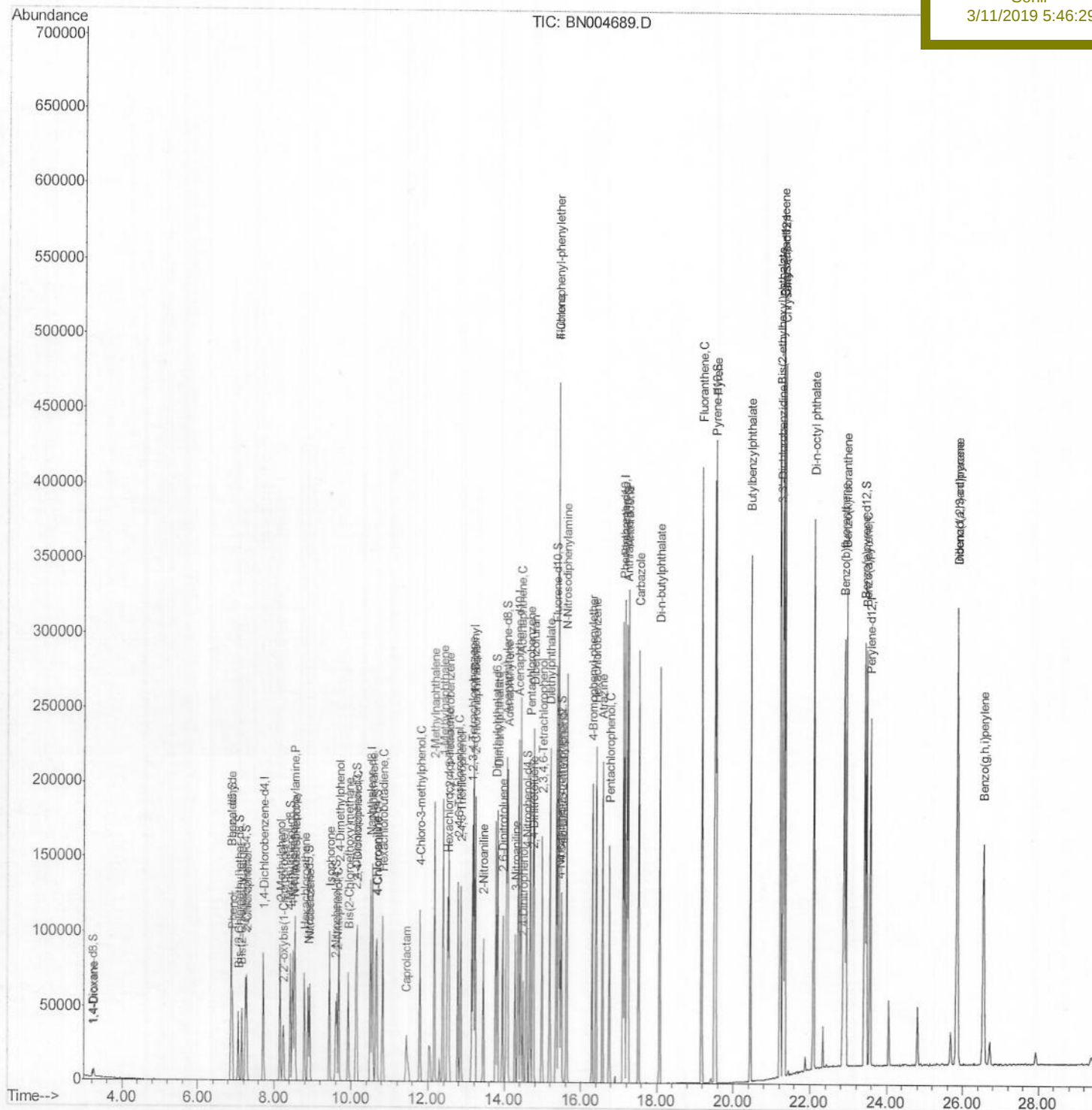
SSTD02027

## Manual Integrations

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3/11/2019 5:46:29 PM



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN030819\

Data File : BN004689.D

Acq On : 09 Mar 2019 00:34

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Sample : SSTDCCC020

Misc :

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Quant Time: Mar 09 01:53:40 2019

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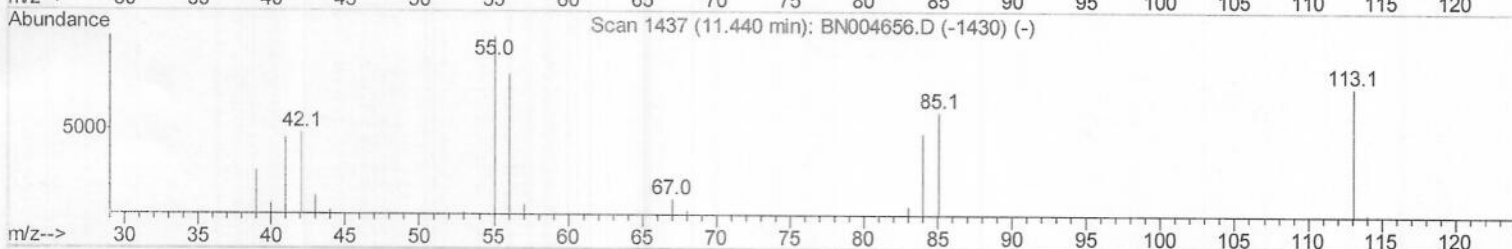
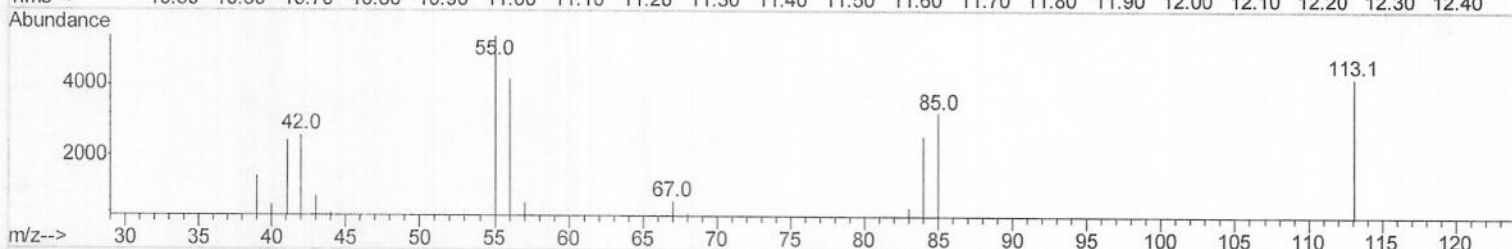
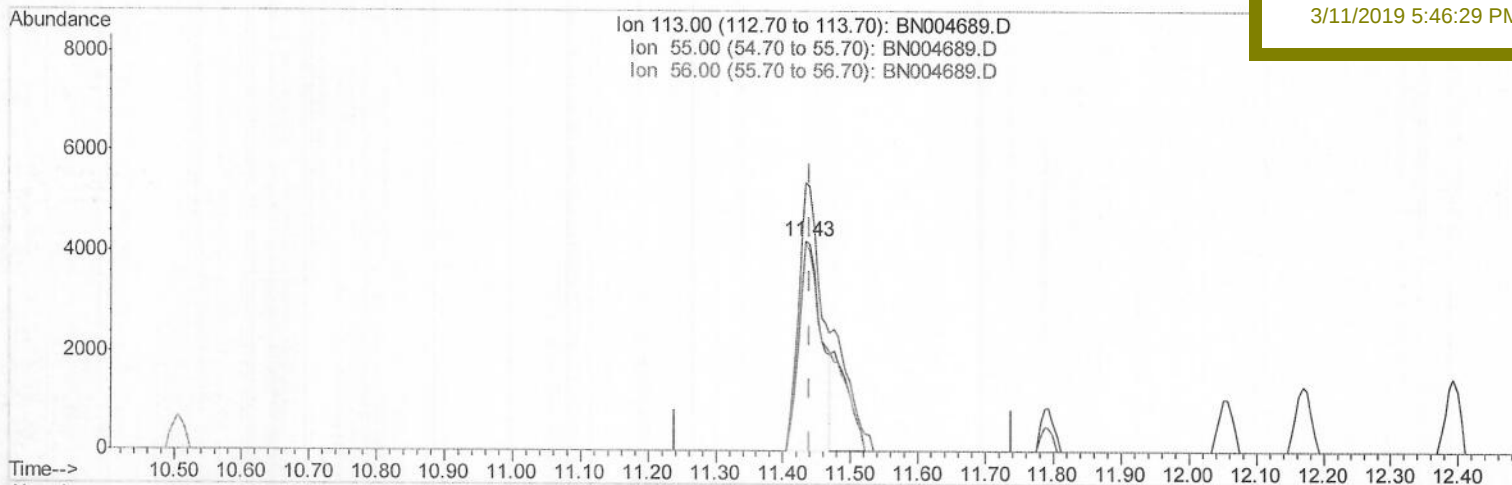
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TIC: BN004689.D

(32) Caprolactam

11.434min (-0.006) 14.71ng/ul

response 9751

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 125.90 | 127.77 |
| 56.00  | 95.10  | 99.29  |
| 0.00   | 0.00   | 0.00   |

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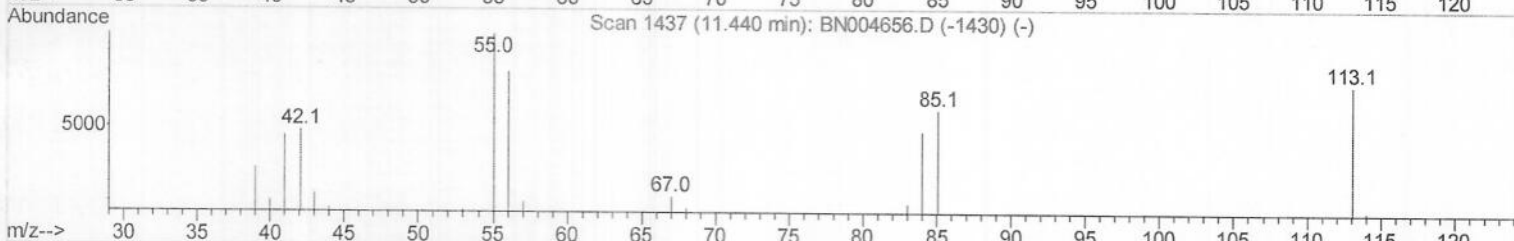
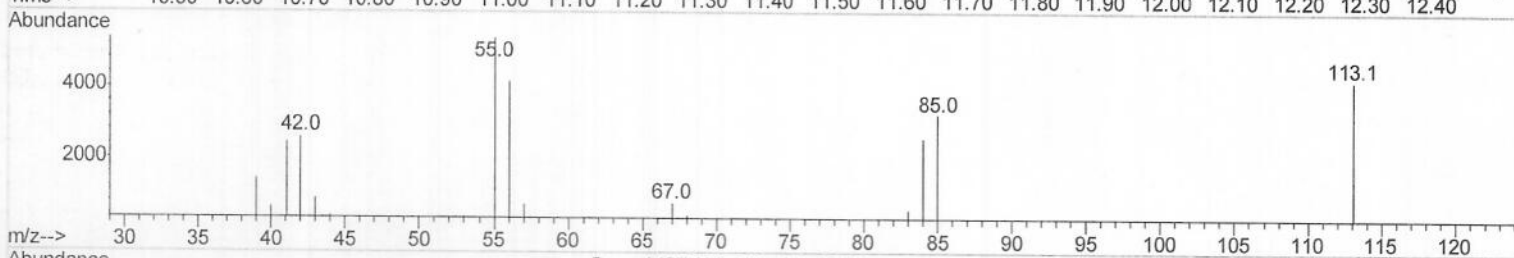
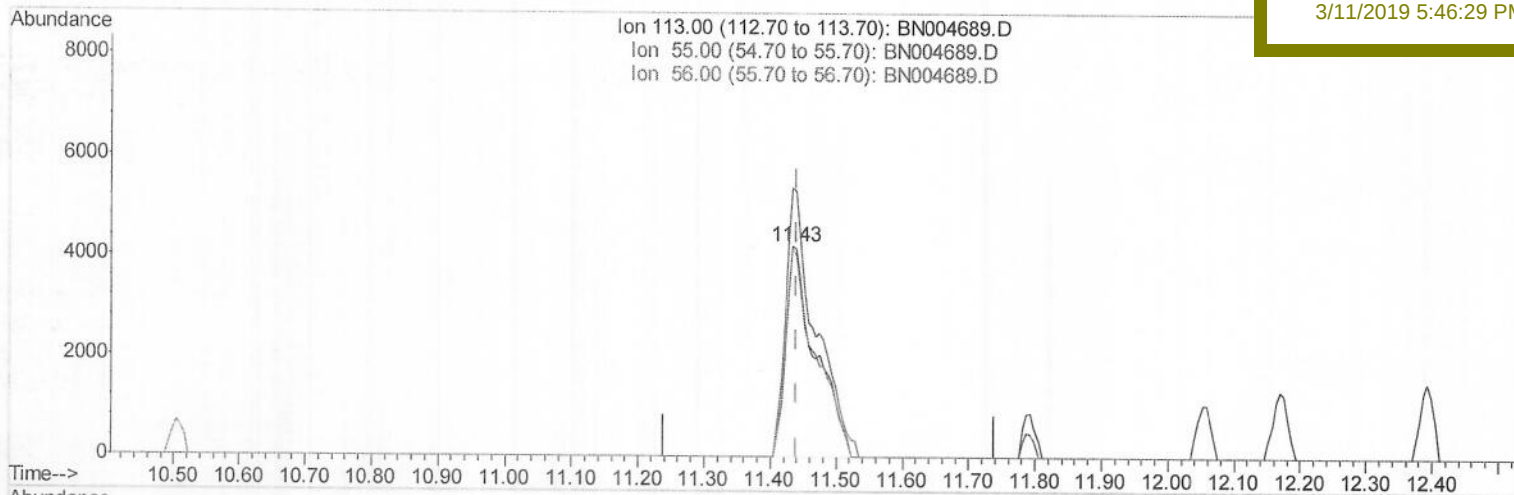
LabSampleId :

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Manual Integrations  
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3/11/2019 5:46:29 PM



TIC: BN004689.D

(32) Caprolactam

11.434min (-0.006) 19.94ng/ul m

response 13215

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 125.90 | 127.77 |
| 56.00  | 95.10  | 99.29  |
| 0.00   | 0.00   | 0.00   |

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03/13/19

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Instrument :  
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Manual Integrations  
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| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.72  | 152  | 25519    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.50 | 136  | 120496   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.36 | 164  | 78051    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.11 | 188  | 187168   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.31 | 240  | 210196   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.57 | 264  | 181871   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.22  | 96  | 3194   | 7.06  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.88  | 99  | 40354  | 19.14 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.06  | 67  | 20221  | 18.38 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.25  | 132 | 35790  | 19.56 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.42  | 113 | 35619  | 19.42 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.88  | 128 | 17979  | 19.57 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.60  | 143 | 20673  | 19.82 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.13 | 165 | 40131  | 20.18 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.65 | 131 | 47297  | 19.83 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.77 | 166 | 127659 | 18.92 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.06 | 160 | 159070 | 19.58 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.57 | 143 | 24083  | 19.23 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.36 | 176 | 112555 | 19.62 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.48 | 200 | 21931  | 16.97 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.21 | 188 | 179249 | 19.66 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.51 | 212 | 206528 | 20.11 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.42 | 264 | 192500 | 19.71 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |        | Ovalue    |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.26  | 88  | 3800   | 6.816  | ng/uL 99  |
| 4) Benzaldehyde                | 6.88  | 77  | 26415  | 19.347 | ng/ul 96  |
| 6) Phenol                      | 6.91  | 94  | 40277  | 18.768 | ng/ul 98  |
| 8) Bis(2-Chloroethyl)ether     | 7.15  | 93  | 29723  | 18.817 | ng/ul 98  |
| 10) 2-Chlorophenol             | 7.29  | 128 | 35761  | 19.370 | ng/ul 99  |
| 11) 2-Methylphenol             | 8.16  | 108 | 33071  | 19.188 | ng/ul 99  |
| 12) 2,2'-oxybis(1-Chloropropan | 8.25  | 45  | 34533  | 18.955 | ng/ul 99  |
| 14) Acetophenone               | 8.55  | 105 | 54357  | 19.765 | ng/ul 99  |
| 15) N-Nitroso-di-n-propylamine | 8.52  | 70  | 25124  | 19.351 | ng/ul 98  |
| 16) 4-Methylphenol             | 8.48  | 108 | 37018  | 19.320 | ng/ul 96  |
| 17) Hexachloroethane           | 8.79  | 117 | 13588  | 19.651 | ng/ul 92  |
| 20) Nitrobenzene               | 8.92  | 77  | 36773  | 19.498 | ng/ul 99  |
| 21) Isophorone                 | 9.44  | 82  | 74796  | 19.142 | ng/ul 99  |
| 23) 2-Nitrophenol              | 9.63  | 139 | 22238  | 19.716 | ng/ul 98  |
| 24) 2,4-Dimethylphenol         | 9.68  | 107 | 43121  | 19.808 | ng/ul 100 |
| 25) Bis(2-Chloroethoxy)methane | 9.92  | 93  | 44661  | 18.776 | ng/ul 100 |
| 27) 2,4-Dichlorophenol         | 10.15 | 162 | 39417  | 20.143 | ng/ul 99  |
| 28) Naphthalene                | 10.56 | 128 | 123500 | 19.671 | ng/ul 99  |
| 30) 4-Chloroaniline            | 10.68 | 127 | 48335  | 19.937 | ng/ul 100 |
| 31) Hexachlorobutadiene        | 10.82 | 225 | 26105  | 20.073 | ng/ul 99  |
| 32) Caprolactam                | 11.43 | 113 | 13215m | 19.936 | ng/ul 99  |
| 33) 4-Chloro-3-methylphenol    | 11.79 | 107 | 43004  | 20.496 | ng/ul 99  |
| 34) 2-Methylnaphthalene        | 12.17 | 142 | 94735  | 20.137 | ng/ul 100 |

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Manual Integrations  
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| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 35) 1-Methylnaphthalene        | 12.39 | 142  | 88514    | 20.036 | ng/ul | 99       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.54 | 216  | 55013    | 20.044 | ng/ul | 97       |
| 38) Hexachlorocyclopentadiene  | 12.51 | 237  | 38017    | 20.107 | ng/ul | 100      |
| 39) 2,4,6-Trichlorophenol      | 12.79 | 196  | 34726    | 19.642 | ng/ul | 99       |
| 40) 2,4,5-Trichlorophenol      | 12.85 | 196  | 38775    | 20.036 | ng/ul | 97       |
| 41) 1,1'-Biphenyl              | 13.19 | 154  | 125619   | 19.131 | ng/ul | 99       |
| 42) 2-Chloronaphthalene        | 13.23 | 162  | 98186    | 19.666 | ng/ul | 99       |
| 43) 2-Nitroaniline             | 13.45 | 65   | 23873    | 19.545 | ng/ul | 99       |
| 45) Dimethylphthalate          | 13.82 | 163  | 125457   | 18.947 | ng/ul | 100      |
| 46) 2,6-Dinitrotoluene         | 13.95 | 165  | 26349    | 19.653 | ng/ul | 94       |
| 48) Acenaphthylene             | 14.09 | 152  | 155030   | 19.529 | ng/ul | 100      |
| 49) 3-Nitroaniline             | 14.28 | 138  | 26495    | 20.381 | ng/ul | 98       |
| 50) Acenaphthene               | 14.43 | 153  | 105977   | 19.530 | ng/ul | 99       |
| 51) 2,4-Dinitrophenol          | 14.49 | 184  | 17106    | 19.686 | ng/ul | 98       |
| 53) 4-Nitrophenol              | 14.58 | 109  | 18512    | 20.341 | ng/ul | 94       |
| 54) Dibenzofuran               | 14.76 | 168  | 153290   | 19.667 | ng/ul | 99       |
| 55) 2,4-Dinitrotoluene         | 14.74 | 165  | 39590    | 19.991 | ng/ul | 98       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.99 | 232  | 35828    | 20.003 | ng/ul | 97       |
| 57) Diethylphthalate           | 15.19 | 149  | 129441   | 19.291 | ng/ul | 99       |
| 59) Fluorene                   | 15.41 | 166  | 123269   | 19.710 | ng/ul | 98       |
| 60) 4-Chlorophenyl-phenylether | 15.41 | 204  | 65460    | 19.715 | ng/ul | 99       |
| 61) 4-Nitroaniline             | 15.45 | 138  | 29772    | 19.511 | ng/ul | 99       |
| 64) 4,6-Dinitro-2-methylphenol | 15.50 | 198  | 23441    | 17.205 | ng/ul | 98       |
| 65) N-Nitrosodiphenylamine     | 15.62 | 169  | 109501   | 19.210 | ng/ul | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.30 | 248  | 44288    | 19.282 | ng/ul | 98       |
| 67) Hexachlorobenzene          | 16.41 | 284  | 52553    | 19.934 | ng/ul | 96       |
| 68) Atrazine                   | 16.57 | 200  | 41871    | 19.231 | ng/ul | 98       |
| 69) Pentachlorophenol          | 16.76 | 266  | 30306    | 18.445 | ng/ul | 98       |
| 70) Phenanthrene               | 17.15 | 178  | 203679   | 19.569 | ng/ul | 100      |
| 72) Anthracene                 | 17.25 | 178  | 209594   | 19.710 | ng/ul | 100      |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.15 | 216  | 54165    | 19.496 | ng/uL | 100      |
| 74) Pentachlorobenzene         | 14.67 | 250  | 58318    | 19.689 | ng/uL | 99       |
| 75) Carbazole                  | 17.52 | 167  | 187371   | 20.289 | ng/ul | 100      |
| 76) Di-n-butylphthalate        | 18.07 | 149  | 220888   | 19.287 | ng/ul | 100      |
| 77) Fluoranthene               | 19.17 | 202  | 251437   | 20.953 | ng/ul | 97       |
| 80) Pyrene                     | 19.54 | 202  | 257565   | 20.212 | ng/ul | 99       |
| 81) Butylbenzylphthalate       | 20.44 | 149  | 107098   | 20.469 | ng/ul | 100      |
| 82) 3,3'-Dichlorobenzidine     | 21.23 | 252  | 98326    | 19.814 | ng/ul | 99       |
| 83) Benzo(a)anthracene         | 21.29 | 228  | 259783   | 19.646 | ng/ul | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.21 | 149  | 157994   | 20.779 | ng/ul | 99       |
| 85) Chrysene                   | 21.35 | 228  | 239994   | 19.545 | ng/ul | 99       |
| 87) Di-n-octyl phthalate       | 22.10 | 149  | 273010   | 23.263 | ng/ul | 100      |
| 88) Benzo(b)fluoranthene       | 22.89 | 252  | 232288   | 20.578 | ng/ul | 99       |
| 89) Benzo(k)fluoranthene       | 22.93 | 252  | 214097   | 19.288 | ng/ul | 99       |
| 91) Benzo(a)pyrene             | 23.47 | 252  | 207930   | 19.580 | ng/ul | 98       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.85 | 276  | 231062   | 19.975 | ng/ul | 98       |
| 93) Dibenzo(a,h)anthracene     | 25.86 | 278  | 195977   | 20.136 | ng/ul | 98       |
| 94) Benzo(a,h,i)perylene       | 26.55 | 276  | 189898   | 20.186 | ng/ul | 96       |

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

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