

(QT Reviewed)

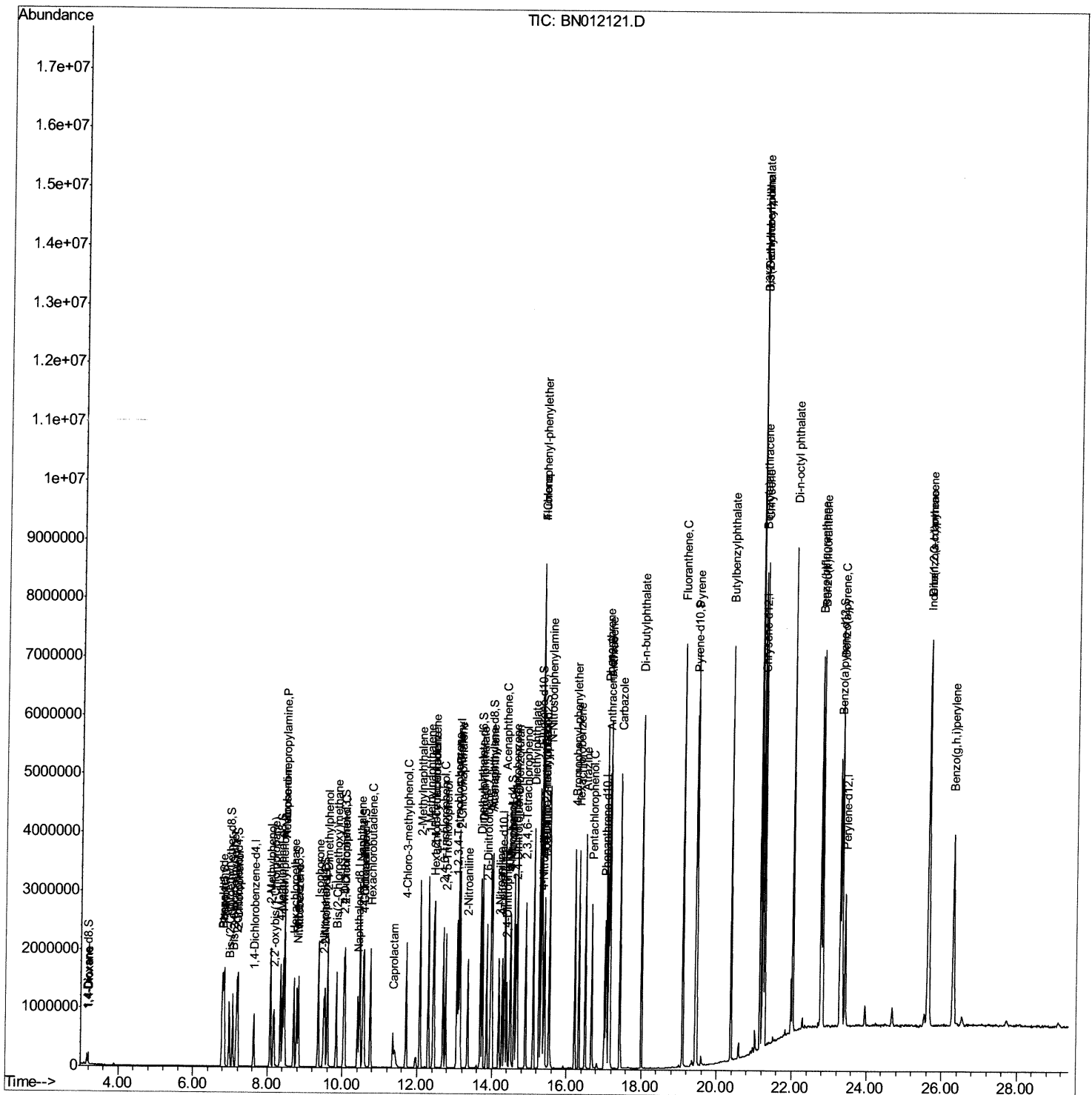
```
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN100920\  
Data File : BN012121.D  
Acq On    : 09 Oct 2020   12:45  
Operator  : CG/JU  
Sample    : SSTD04056  
Misc      :  
ALS Vial  : 5      Sample Multiplier: 1
```

**Instrument :**  
BNA\_N  
**ClientSampleId :**  
SSTD04056

## Manual Integrations

mohammad  
10/12/2020 11:35:15 AM

Quant Time: Oct 09 13:21:24 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN100920.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Oct 09 12:44:31 2020  
Response via : Initial Calibration



# Quantitation Report (Qedit)

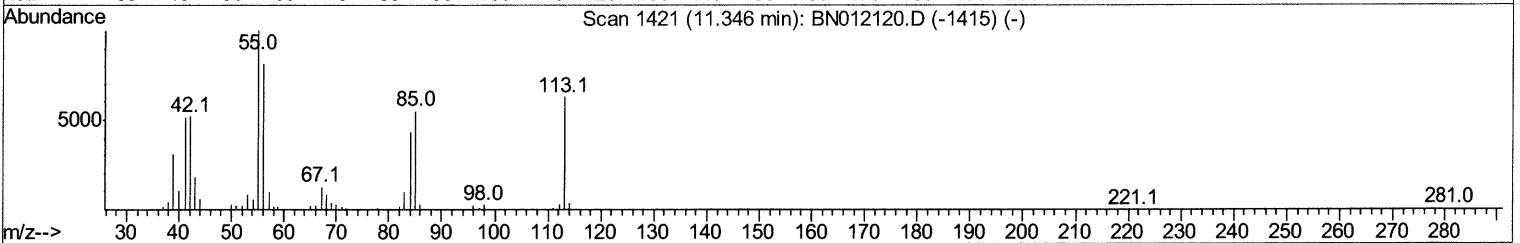
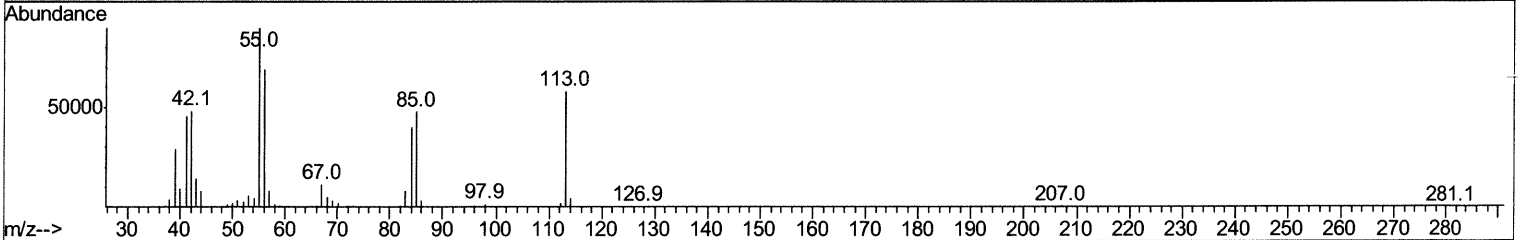
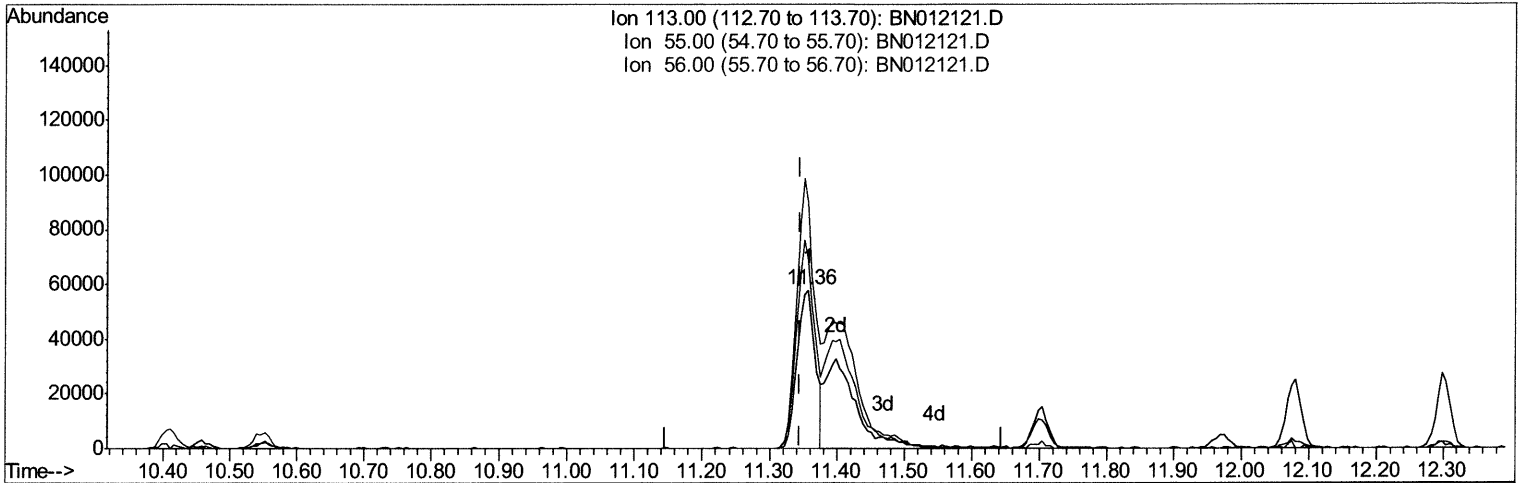
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN100920\  
 Data File : BN012121.D  
 Acq On : 09 Oct 2020 12:45  
 Operator : CG/JU  
 Sample : SSTD04056  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

**Instrument :**  
 BNA\_N  
**ClientSampleId :**  
 SSTD04056

**Manual Integrations**  
**APPROVED**

mohammad  
 10/12/2020 11:35:15 AM

Quant Time: Oct 09 13:20:30 2020  
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TIC: BN012121.D

(32) Caprolactam

11.357min (+0.012) 25.58ng/ul

response 108352

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 158.80 | 154.63 |
| 56.00  | 129.40 | 118.84 |
| 0.00   | 0.00   | 0.00   |

```
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Operator  : CG/JU  
Sample    : SSTD04056  
Misc      :  
ALS Vial  : 5      Sample Multiplier: 1
```

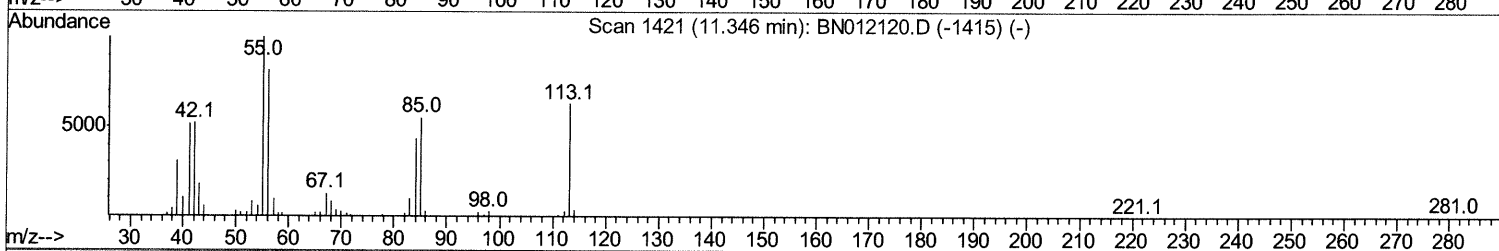
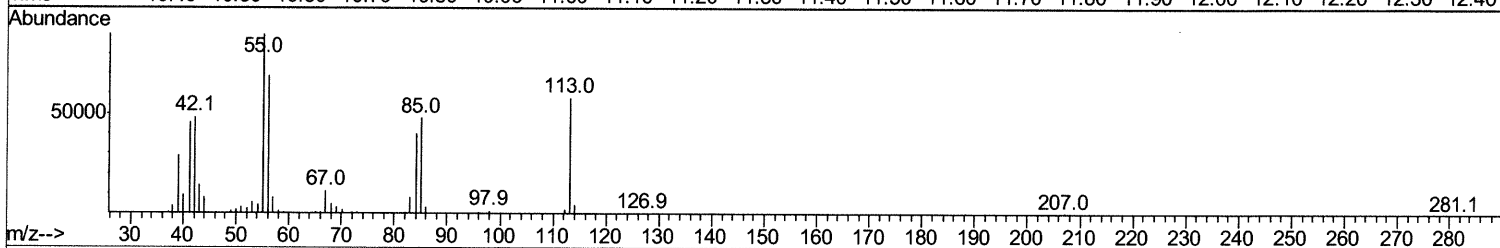
mohammad  
10/12/2020 11:35:15 AM

Abundance

Ion 113.00 (112.70 to 113.70): BN012121.D  
 Ion 55.00 (54.70 to 55.70): BN012121.D  
 Ion 56.00 (55.70 to 56.70): BN012121.D

1d 3d 4d

Time-->



TIC: BN012121.D

|        |        |        |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 158.80 | 154.63 |
| 56.00  | 129.40 | 118.84 |
| 0.00   | 0.00   | 0.00   |

<sup>m</sup> Ju 10/15/20

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100920\  
 Data File : BN012121.D  
 Acq On : 09 Oct 2020 12:45  
 Operator : CG/JU  
 Sample : SST04056  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST04056

Manual Integrations  
 APPROVED

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Quant Time: Oct 09 13:21:24 2020  
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 Quant Title : SVOA CALIBRATION  
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 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.63  | 152  | 226888   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.41 | 136  | 920956   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.27 | 164  | 600884   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.03 | 188  | 1391385  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.23 | 240  | 1426419  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.43 | 264  | 1460335  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.16  | 96  | 78415   | 15.04 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.81  | 99  | 750317  | 39.62 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.98  | 67  | 463619  | 39.85 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.17  | 132 | 613731  | 40.08 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.34  | 113 | 618620  | 41.50 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.79  | 128 | 298558  | 40.83 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.50  | 143 | 348015  | 44.00 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.03 | 165 | 610966  | 40.28 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.55 | 131 | 819982  | 45.86 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.69 | 166 | 1977479 | 41.95 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.97 | 160 | 2350963 | 41.79 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.49 | 143 | 386322  | 44.51 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.27 | 176 | 1703625 | 41.42 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.40 | 200 | 368061  | 41.45 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.13 | 188 | 2656710 | 39.38 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.43 | 212 | 2985354 | 39.79 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.29 | 264 | 3277193 | 39.99 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |              | Ovalue |
|--------------------------------|-------|-----|---------|--------------|--------|
| 2) 1,4-Dioxane                 | 3.20  | 88  | 88404   | 13.503 ng/uL | 92     |
| 4) Benzaldehyde                | 6.79  | 77  | 505928  | 46.413 ng/ul | 97     |
| 6) Phenol                      | 6.84  | 94  | 777561  | 39.476 ng/ul | 97     |
| 8) Bis(2-Chloroethyl)ether     | 7.07  | 93  | 632625  | 39.532 ng/ul | 97     |
| 10) 2-Chlorophenol             | 7.20  | 128 | 639035  | 40.349 ng/ul | 98     |
| 11) 2-Methylphenol             | 8.08  | 108 | 601721  | 41.128 ng/ul | 98     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.16  | 45  | 823989  | 39.926 ng/ul | 100    |
| 14) Acetophenone               | 8.45  | 105 | 955710  | 40.229 ng/ul | 96     |
| 15) N-Nitroso-di-n-propylamine | 8.45  | 70  | 511104  | 42.873 ng/ul | 99     |
| 16) 4-Methylphenol             | 8.40  | 108 | 653477  | 41.653 ng/ul | 95     |
| 17) Hexachloroethane           | 8.70  | 117 | 276386  | 38.356 ng/ul | 94     |
| 20) Nitrobenzene               | 8.83  | 77  | 794853  | 39.324 ng/ul | 95     |
| 21) Isophorone                 | 9.35  | 82  | 1439955 | 42.502 ng/ul | 98     |
| 23) 2-Nitrophenol              | 9.53  | 139 | 368488  | 42.123 ng/ul | 98     |
| 24) 2,4-Dimethylphenol         | 9.60  | 107 | 756169  | 39.836 ng/ul | 95     |
| 25) Bis(2-Chloroethoxy)methane | 9.83  | 93  | 861694  | 40.530 ng/ul | 99     |
| 27) 2,4-Dichlorophenol         | 10.06 | 162 | 622811  | 40.550 ng/ul | 99     |
| 28) Naphthalene                | 10.46 | 128 | 2100377 | 39.848 ng/ul | 99     |
| 30) 4-Chloroaniline            | 10.57 | 127 | 790853  | 43.916 ng/ul | 98     |
| 31) Hexachlorobutadiene        | 10.75 | 225 | 394254  | 37.913 ng/ul | 98     |
| 32) Caprolactam                | 11.36 | 113 | 202715m | 47.864 ng/ul | 98     |
| 33) 4-Chloro-3-methylphenol    | 11.70 | 107 | 725367  | 43.458 ng/ul | 98     |
| 34) 2-Methylnaphthalene        | 12.08 | 142 | 1496958 | 41.324 ng/ul | 100    |

Jul 10/15/20

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 Quant Title : SVOA CALIBRATION  
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| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.30 | 142  | 1438229  | 40.177 | ng/uL  | 99       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.45 | 216  | 713045   | 36.631 | ng/uL  | 91       |
| 38) Hexachlorocyclopentadiene  | 12.43 | 237  | 490308   | 35.624 | ng/uL  | 98       |
| 39) 2,4,6-Trichlorophenol      | 12.70 | 196  | 504643   | 41.350 | ng/uL  | 96       |
| 40) 2,4,5-Trichlorophenol      | 12.77 | 196  | 540799   | 41.078 | ng/uL  | 93       |
| 41) 1,1'-Biphenyl              | 13.10 | 154  | 1900056  | 38.174 | ng/uL  | 100      |
| 42) 2-Chloronaphthalene        | 13.15 | 162  | 1539392  | 38.835 | ng/uL  | 98       |
| 43) 2-Nitroaniline             | 13.35 | 65   | 497230   | 43.772 | ng/uL  | 97       |
| 45) Dimethylphthalate          | 13.74 | 163  | 2041632  | 42.035 | ng/uL  | 99       |
| 46) 2,6-Dinitrotoluene         | 13.86 | 165  | 436711   | 46.065 | ng/uL  | 95       |
| 48) Acenaphthylene             | 14.00 | 152  | 2385617  | 40.355 | ng/uL  | 97       |
| 49) 3-Nitroaniline             | 14.19 | 138  | 416350   | 48.131 | ng/uL  | 97       |
| 50) Acenaphthene               | 14.34 | 153  | 1693507  | 39.859 | ng/uL  | 97       |
| 51) 2,4-Dinitrophenol          | 14.39 | 184  | 272959   | 47.227 | ng/uL  | 93       |
| 53) 4-Nitrophenol              | 14.50 | 109  | 342261   | 42.322 | ng/uL  | 97       |
| 54) Dibenzofuran               | 14.68 | 168  | 2373522  | 40.947 | ng/uL  | 99       |
| 55) 2,4-Dinitrotoluene         | 14.65 | 165  | 626992   | 45.560 | ng/uL  | 100      |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.90 | 232  | 487223   | 44.338 | ng/uL# | 97       |
| 57) Diethylphthalate           | 15.12 | 149  | 2162277  | 43.257 | ng/uL  | 99       |
| 59) Fluorene                   | 15.33 | 166  | 2007767  | 41.317 | ng/uL  | 98       |
| 60) 4-Chlorophenyl-phenylether | 15.33 | 204  | 961503   | 40.539 | ng/uL  | 99       |
| 61) 4-Nitroaniline             | 15.36 | 138  | 462187   | 49.447 | ng/uL  | 96       |
| 64) 4,6-Dinitro-2-methylphenol | 15.42 | 198  | 397524   | 41.157 | ng/uL# | 96       |
| 65) N-Nitrosodiphenylamine     | 15.55 | 169  | 1686815  | 38.012 | ng/uL  | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.22 | 248  | 596863   | 37.508 | ng/uL  | 92       |
| 67) Hexachlorobenzene          | 16.33 | 284  | 728856   | 39.253 | ng/uL  | 98       |
| 68) Atrazine                   | 16.50 | 200  | 655781   | 40.390 | ng/uL  | 96       |
| 69) Pentachlorophenol          | 16.68 | 266  | 447183   | 40.620 | ng/uL  | 95       |
| 70) Phenanthrene               | 17.07 | 178  | 3260256  | 38.462 | ng/uL  | 96       |
| 72) Anthracene                 | 17.16 | 178  | 3405820  | 39.791 | ng/uL  | 98       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.06 | 216  | 722289   | 35.710 | ng/uL  | 97       |
| 74) Pentachlorobenzene         | 14.60 | 250  | 762046   | 36.731 | ng/uL  | 96       |
| 75) Carbazole                  | 17.43 | 167  | 2949814  | 39.691 | ng/uL  | 99       |
| 76) Di-n-butylphthalate        | 18.00 | 149  | 3918166  | 41.517 | ng/uL  | 99       |
| 77) Fluoranthene               | 19.09 | 202  | 4015730  | 42.178 | ng/uL  | 100      |
| 80) Perylene                   | 19.46 | 202  | 4044472  | 39.308 | ng/uL  | 99       |
| 81) Butylbenzylphthalate       | 20.37 | 149  | 1849531  | 43.724 | ng/uL  | 99       |
| 82) 3,3'-Dichlorobenzidine     | 21.16 | 252  | 1469141  | 44.271 | ng/uL  | 95       |
| 83) Benzo(a)anthracene         | 21.22 | 228  | 3797852  | 39.522 | ng/uL  | 97       |
| 84) Bis(2-ethylhexyl)phthalate | 21.16 | 149  | 2870268  | 44.213 | ng/uL# | 96       |
| 85) Chrysene                   | 21.27 | 228  | 3686859  | 39.027 | ng/uL  | 99       |
| 87) Di-n-octyl phthalate       | 22.03 | 149  | 4776049  | 44.034 | ng/uL  | 100      |
| 88) Benzo(b)fluoranthene       | 22.77 | 252  | 3812895  | 37.271 | ng/uL  | 97       |
| 89) Benzo(k)fluoranthene       | 22.82 | 252  | 3876571  | 38.469 | ng/uL  | 98       |
| 91) Benzo(a)pyrene             | 23.34 | 252  | 3559818  | 38.360 | ng/uL  | 98       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.63 | 276  | 4473839  | 37.755 | ng/uL  | 98       |
| 93) Dibenzo(a,h)anthracene     | 25.64 | 278  | 3780222  | 38.723 | ng/uL  | 97       |
| 94) Benzo(a,h,i)perylene       | 26.30 | 276  | 3798756  | 37.643 | ng/uL  | 99       |

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 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD04056

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 QLast Update : Fri Oct 09 12:44:31 2020  
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| Internal Standards                                                         | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------------------------------------------------------|------|------|----------|------|-------|----------|
| -----                                                                      |      |      |          |      |       |          |
| (#) = qualifier out of range (m) = manual integration (+) = signals summed |      |      |          |      |       |          |