

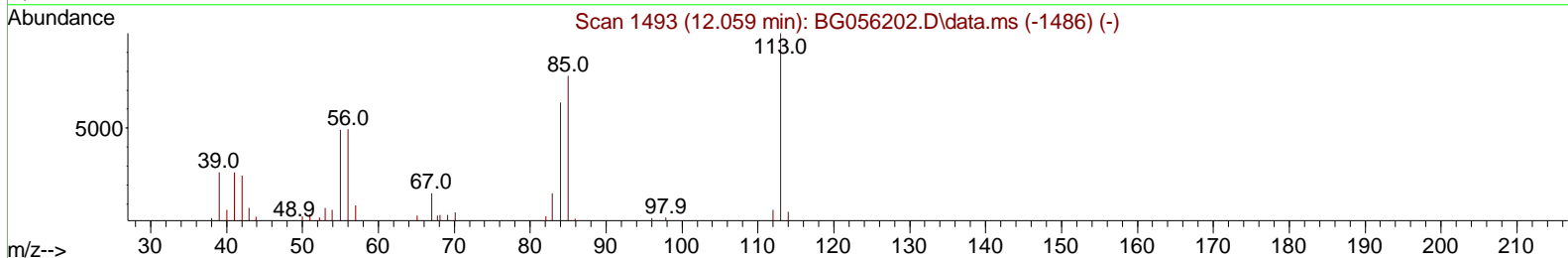
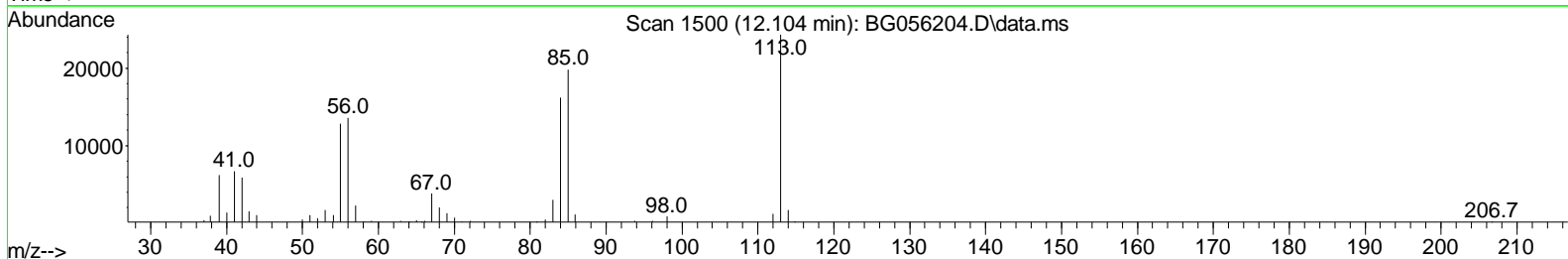
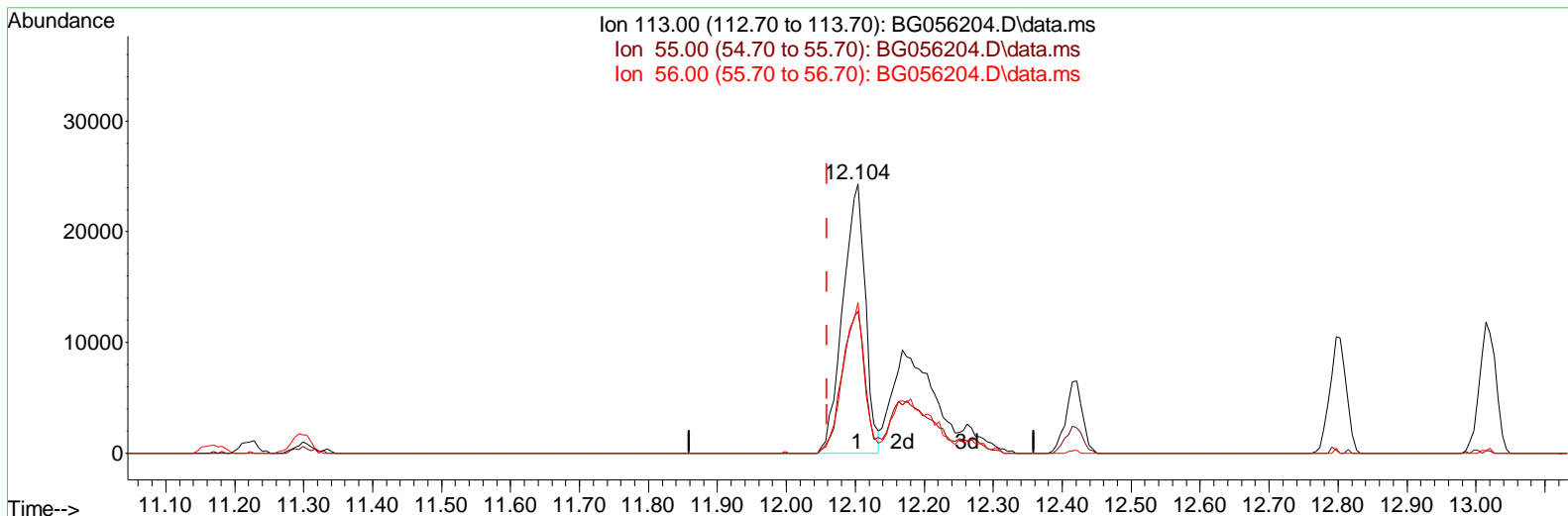
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG010423\  
 Data File : BG056204.D  
 Acq On : 4 Jan 2023 17:47  
 Operator : CG/JU  
 Sample : SSTD08051  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD080404

Manual Integrations APPROVED

Reviewed By : Jagrut Upadhyay 01/05/2023  
 Supervised By : mohammad ahmed 01/05/2023

Quant Time: Jan 05 00:20:18 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010423.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jan 05 00:14:25 2023  
 Response via : Initial Calibration



TIC: BG056204.D\data.ms

(34) Caprolactam

12.104min (+ 0.045) 39.05 ng/ul

response 55254

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 49.10  | 52.80  |
| 56.00  | 49.40  | 56.05  |
| 0.00   | 0.00   | 0.00   |

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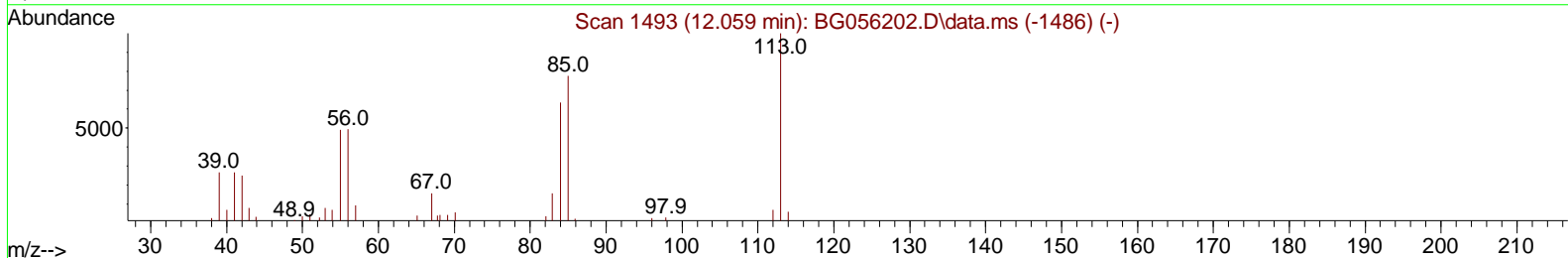
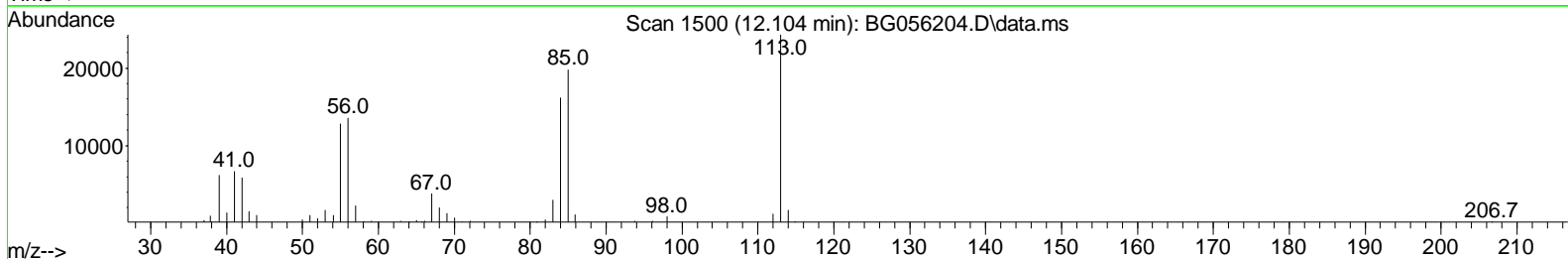
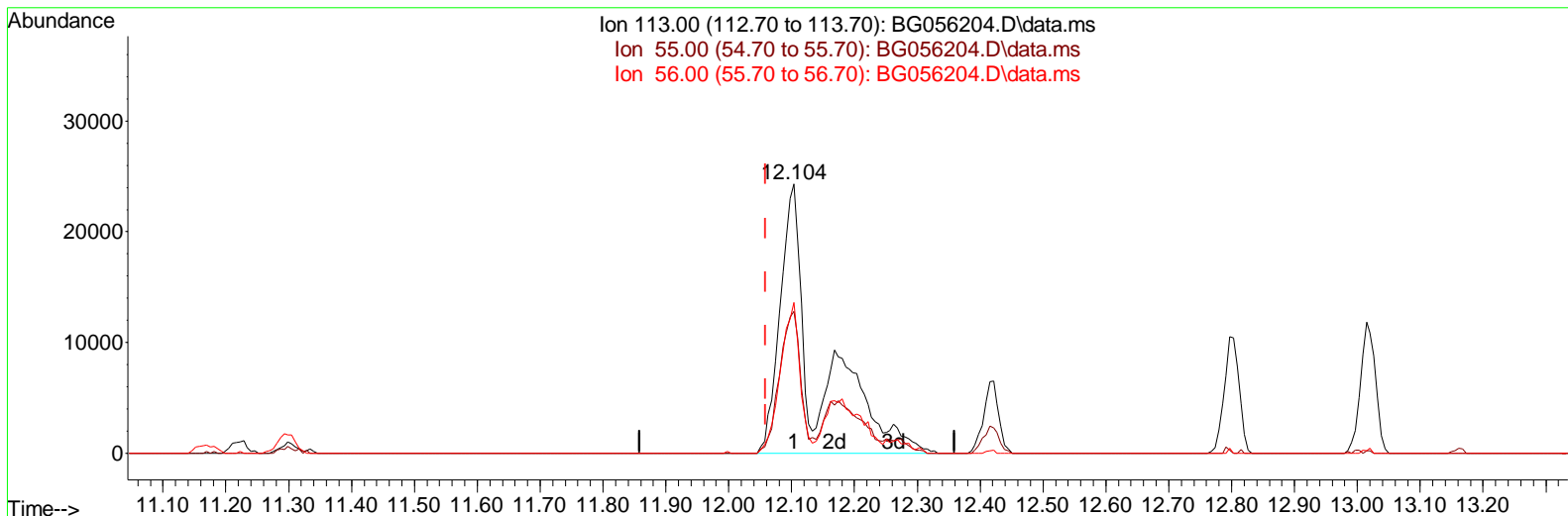
**ClientSampleId :**

SSTD080404

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TIC: BG056204.D\data.ms

**(34) Caprolactam**

12.104min (+ 0.045) 69.92 ng/ul m

response 98926

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 49.10  | 52.80  |
| 56.00  | 49.40  | 56.05  |
| 0.00   | 0.00   | 0.00   |

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**Instrument :**  
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**ClientSampleId :**  
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 Quant Title : SVOA CALIBRATION  
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 Response via : Initial Calibration

| Compound                           | R. T.  | QI on | Response | Conc    | Units   | Dev(Min) |
|------------------------------------|--------|-------|----------|---------|---------|----------|
| <b>Internal Standards</b>          |        |       |          |         |         |          |
| 1) 1,4-Dichlorobenzene-d4          | 8.332  | 152   | 54294    | 20.000  | ng/ul   | 0.00     |
| 20) Naphthalene-d8                 | 11.170 | 136   | 219125   | 20.000  | ng/ul   | 0.00     |
| 38) Acenaphthene-d10               | 14.947 | 164   | 173322   | 20.000  | ng/ul   | 0.00     |
| 64) Phenanthrene-d10               | 17.685 | 188   | 417979   | 20.000  | ng/ul   | 0.00     |
| 79) Chrysene-d12                   | 21.986 | 240   | 398609   | 20.000  | ng/ul   | 0.00     |
| 88) Perylene-d12                   | 25.447 | 264   | 437779   | 20.000  | ng/ul   | 0.02     |
| <b>System Monitoring Compounds</b> |        |       |          |         |         |          |
| 3) 1,4-Dioxane-d8                  | 3.631  | 96    | 37686    | 29.055  | ng/uL   | 0.00     |
| 4) Pyridine-d5                     | 4.066  | 84    | 302740   | 71.587  | ng/ul   | 0.00     |
| 7) Phenol-d5                       | 7.468  | 99    | 385267   | 78.227  | ng/ul   | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...       | 7.644  | 67    | 154672   | 79.683  | ng/ul   | 0.00     |
| 11) 2-Chlorophenol-d4              | 7.862  | 132   | 263133   | 90.399  | ng/ul   | 0.00     |
| 15) 4-Methylphenol-d8              | 9.037  | 113   | 302273   | 81.551  | ng/ul   | 0.00     |
| 21) Nitrobenzene-d5                | 9.513  | 128   | 138658   | 86.879  | ng/ul   | 0.00     |
| 24) 2-Nitrophenol-d4               | 10.241 | 143   | 180703   | 86.660  | ng/ul   | 0.00     |
| 28) 2,4-Dichlorophenol-d3          | 10.782 | 165   | 360411   | 77.781  | ng/ul   | 0.00     |
| 31) 4-Chloroaniline-d4             | 11.299 | 131   | 388622   | 74.149  | ng/ul   | 0.00     |
| 46) Dimethylphthalate-d6           | 14.342 | 166   | 1055854  | 78.711  | ng/ul   | 0.00     |
| 49) Acenaphthylene-d8              | 14.648 | 160   | 1185657  | 78.265  | ng/ul   | 0.00     |
| 54) 4-Nitrophenol-d4               | 15.130 | 143   | 165454   | 81.938  | ng/ul   | 0.02     |
| 60) Fluorene-d10                   | 15.935 | 176   | 995204   | 75.160  | ng/ul   | 0.00     |
| 65) 4,6-Dinitro-2-methylph...      | 16.046 | 200   | 244929   | 91.156  | ng/ul   | 0.02     |
| 73) Anthracene-d10                 | 17.785 | 188   | 1488383  | 77.057  | ng/ul   | 0.00     |
| 81) Pyrene-d10                     | 20.047 | 212   | 1749706  | 79.481  | ng/ul   | 0.00     |
| 92) Benzo(a)pyrene-d12             | 25.218 | 264   | 1773295  | 79.694  | ng/ul   | 0.02     |
| <b>Target Compounds</b>            |        |       |          |         |         |          |
| 2) 1,4-Dioxane                     | 3.672  | 88    | 38415    | 26.477  | ng/uL   | 98       |
| 5) Pyridine                        | 4.090  | 79    | 297368   | 73.114  | ng/ul   | 97       |
| 6) Benzaldehyde                    | 7.462  | 77    | 131357   | 72.650  | ng/ul   | 97       |
| 8) Phenol                          | 7.497  | 94    | 378262   | 77.174  | ng/ul   | 96       |
| 10) Bis(2-Chloroethyl)ether        | 7.744  | 93    | 256633   | 79.543  | ng/ul   | 98       |
| 12) 2-Chlorophenol                 | 7.897  | 128   | 259435   | 88.736  | ng/ul   | 96       |
| 13) 2-Methylphenol                 | 8.766  | 108   | 292934   | 79.598  | ng/ul   | 95       |
| 14) 2,2'-oxybis(1-Chloropr...      | 8.855  | 45    | 47571    | 114.793 | ng/ul # | 81       |
| 16) Acetophenone                   | 9.172  | 105   | 461927   | 75.236  | ng/ul   | 96       |
| 17) N-Nitrosodipropylamine         | 9.160  | 70    | 205634   | 79.216  | ng/ul   | 99       |
| 18) 4-Methylphenol                 | 9.107  | 108   | 316592   | 79.418  | ng/ul   | 99       |
| 19) Hexachloroethane               | 9.436  | 117   | 99182    | 84.374  | ng/ul   | 95       |
| 22) Nitrobenzene                   | 9.560  | 77    | 313456   | 79.151  | ng/ul   | 96       |
| 23) Isophorone                     | 10.088 | 82    | 639022   | 70.574  | ng/ul   | 99       |
| 25) 2-Nitrophenol                  | 10.276 | 139   | 171436   | 85.170  | ng/ul   | 99       |
| 26) 2,4-Dimethylphenol             | 10.318 | 107   | 354121   | 74.349  | ng/ul   | 96       |
| 27) Bis(2-Chloroethoxy)meth...     | 10.553 | 93    | 354377   | 75.521  | ng/ul   | 96       |
| 29) 2,4-Dichlorophenol             | 10.811 | 162   | 340476   | 76.933  | ng/ul   | 97       |
| 30) Naphthalene                    | 11.222 | 128   | 897333   | 78.593  | ng/ul   | 99       |
| 32) 4-Chloroaniline                | 11.322 | 127   | 396522   | 75.079  | ng/ul   | 96       |
| 33) Hexachlorobutadiene            | 11.493 | 225   | 288181   | 69.486  | ng/ul   | 99       |
| 34) Caprolactam                    | 12.104 | 113   | 98926m   | 69.917  | ng/ul   | 99       |
| 35) 4-Chloro-3-methylphenol        | 12.415 | 107   | 317249   | 74.617  | ng/ul   | 95       |

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 Sample : SSTD08051  
 Misc :  
 ALS Vial : 7 Sample Multi plier: 1

**Instrument :**  
 BNA\_G  
**ClientSampleId :**  
 SSTD080404

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 01/05/2023  
 Supervised By :mohammad ahmed 01/05/2023

Quant Time: Jan 05 00:20:18 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010423.M  
 Quant Title : SVOA CALI BRATI ON  
 QLast Update : Thu Jan 05 00:14:25 2023  
 Response via : Ini tial Cal i brati on

| Compound                            | R. T.  | QI on | Response | Conc    | Units   | Dev(Mi n) |
|-------------------------------------|--------|-------|----------|---------|---------|-----------|
| 36) 2-Methyl naphthal ene           | 12.803 | 142   | 680177   | 74.305  | ng/ul   | 99        |
| 37) 1-Methyl naphthal ene           | 13.014 | 142   | 681535   | 73.908  | ng/ul   | 94        |
| 39) 1, 2, 4, 5-Tetrachl oroben. . . | 13.161 | 216   | 545447   | 76.307  | ng/ul   | 100       |
| 40) Hexachl orocycl opentadi ene    | 13.138 | 237   | 389035   | 83.939  | ng/ul   | 96        |
| 41) 2, 4, 6-Tri chl orophenol       | 13.390 | 196   | 352493   | 83.667  | ng/ul   | 98        |
| 42) 2, 4, 5-Tri chl orophenol       | 13.467 | 196   | 378179   | 82.788  | ng/ul   | 96        |
| 43) 1, 1' -Bi phenyl                | 13.790 | 154   | 980608   | 80.028  | ng/ul   | 99        |
| 44) 2-Chl oronaphthal ene           | 13.837 | 162   | 818559   | 80.596  | ng/ul   | 96        |
| 45) 2-Ni troani li ne               | 14.031 | 65    | 156761   | 111.147 | ng/ul   | 91        |
| 47) Di methyl phthal ate            | 14.389 | 163   | 1014451  | 76.263  | ng/ul   | 100       |
| 48) 2, 6-Di ni trotol uene          | 14.519 | 165   | 226349   | 86.209  | ng/ul   | 95        |
| 50) Acenaphthyl ene                 | 14.677 | 152   | 1181561  | 78.483  | ng/ul   | 97        |
| 51) 3-Ni troani li ne               | 14.848 | 138   | 168559   | 85.467  | ng/ul   | 98        |
| 52) Acenaphthene                    | 15.012 | 153   | 835444   | 78.993  | ng/ul   | 98        |
| 53) 2, 4-Di ni trophenol            | 15.047 | 184   | 171592   | 83.963  | ng/ul # | 91        |
| 55) 4-Ni trophenol                  | 15.147 | 109   | 161632   | 82.276  | ng/ul   | 92        |
| 56) Di benzofuran                   | 15.347 | 168   | 1243997  | 75.147  | ng/ul   | 97        |
| 57) 2, 4-Di ni trotol uene          | 15.300 | 165   | 310238   | 85.207  | ng/ul # | 97        |
| 58) 2, 3, 4, 6-Tetrachl orophenol   | 15.564 | 232   | 358001   | 79.285  | ng/ul # | 98        |
| 59) Di ethyl phthal ate             | 15.741 | 149   | 935666   | 78.733  | ng/ul   | 97        |
| 61) Fl uorene                       | 15.993 | 166   | 1000741  | 73.592  | ng/ul   | 98        |
| 62) 4-Chl orophenyl -phenyl e. . .  | 15.970 | 204   | 632419   | 71.588  | ng/ul   | 94        |
| 63) 4-Ni troani li ne               | 16.011 | 138   | 156332   | 83.072  | ng/ul   | 87        |
| 66) 4, 6-Di ni tro-2-methyl ph. . . | 16.064 | 198   | 230082   | 89.314  | ng/ul   | 97        |
| 67) N-Ni trosodi phenyl ami ne      | 16.181 | 169   | 900752   | 85.054  | ng/ul   | 99        |
| 68) 4-Bromophenyl -phenyl ether     | 16.863 | 248   | 437744   | 82.607  | ng/ul   | 95        |
| 69) Hexachl orobenzene              | 16.992 | 284   | 434587   | 84.736  | ng/ul   | 98        |
| 70) Atrazi ne                       | 17.121 | 200   | 397951   | 78.810  | ng/ul   | 100       |
| 71) Pentachl orophenol              | 17.327 | 266   | 276049   | 84.659  | ng/ul   | 98        |
| 72) Phenanthrene                    | 17.727 | 178   | 1657315  | 78.104  | ng/ul   | 97        |
| 74) Anthracene                      | 17.821 | 178   | 1647791  | 75.858  | ng/ul   | 97        |
| 75) 1, 2, 3, 4-Tetrachl oroben. . . | 13.761 | 216   | 553348   | 87.277  | ng/uL   | 98        |
| 76) Pentachl orobenzene             | 15.265 | 250   | 541735   | 86.558  | ng/uL   | 95        |
| 77) Carbazol e                      | 18.085 | 167   | 1391519  | 78.354  | ng/ul   | 99        |
| 78) Di -n-butyl phthal ate          | 18.608 | 149   | 1451700  | 89.470  | ng/ul   | 97        |
| 80) Fl uoranthene                   | 19.718 | 202   | 2057898  | 79.498  | ng/ul   | 99        |
| 82) Pyrene                          | 20.083 | 202   | 1998508  | 76.531  | ng/ul   | 99        |
| 83) Butyl benzyl phthal ate         | 20.935 | 149   | 639624   | 115.098 | ng/ul   | 97        |
| 84) 3, 3' -Di chl orobenzi di ne    | 21.863 | 252   | 706445   | 78.880  | ng/ul   | 99        |
| 85) Benzo(a)anthracene              | 21.963 | 228   | 2111120  | 77.595  | ng/ul   | 97        |
| 86) Bi s(2-ethyl hexyl )phtha. . .  | 21.828 | 149   | 930014   | 109.244 | ng/ul   | 99        |
| 87) Chrysene                        | 22.039 | 228   | 1973325  | 78.153  | ng/ul   | 97        |
| 89) Di -n-octyl phthal ate          | 23.120 | 149   | 1562034  | 96.419  | ng/ul   | 100       |
| 90) Benzo(b)fl uoranthene           | 24.348 | 252   | 2265400  | 78.387  | ng/ul   | 99        |
| 91) Benzo(k)fl uoranthene           | 24.425 | 252   | 2200840  | 78.040  | ng/ul   | 100       |
| 93) Benzo(a)pyrene                  | 25.300 | 252   | 1971550  | 78.289  | ng/ul   | 98        |
| 94) I ndeno(1, 2, 3-cd)pyrene       | 29.454 | 276   | 2651133  | 79.928  | ng/ul   | 99        |
| 95) Di benzo(a, h)anthracene        | 29.530 | 278   | 2198974  | 79.262  | ng/ul   | 100       |
| 96) Benzo(g, h, i )peryl ene        | 30.723 | 276   | 2136529  | 80.151  | ng/ul   | 99        |

(#) = qual i fier out of range (m) = manual i ntegrati on (+) = signal s summed

**Instrument :**

BNA\_G

**ClientSampleId :**

SSTD080404

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