

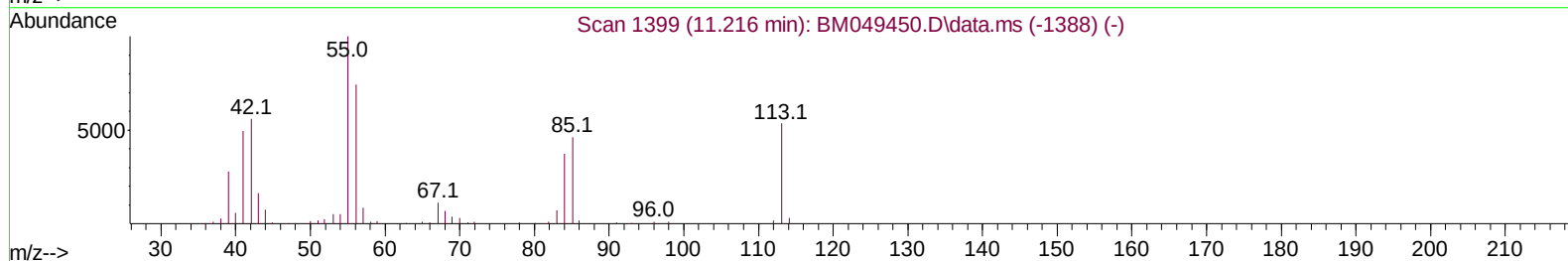
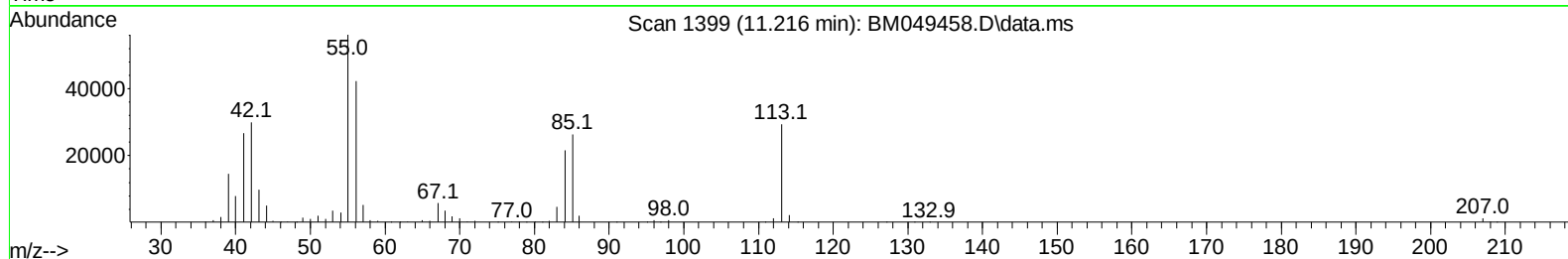
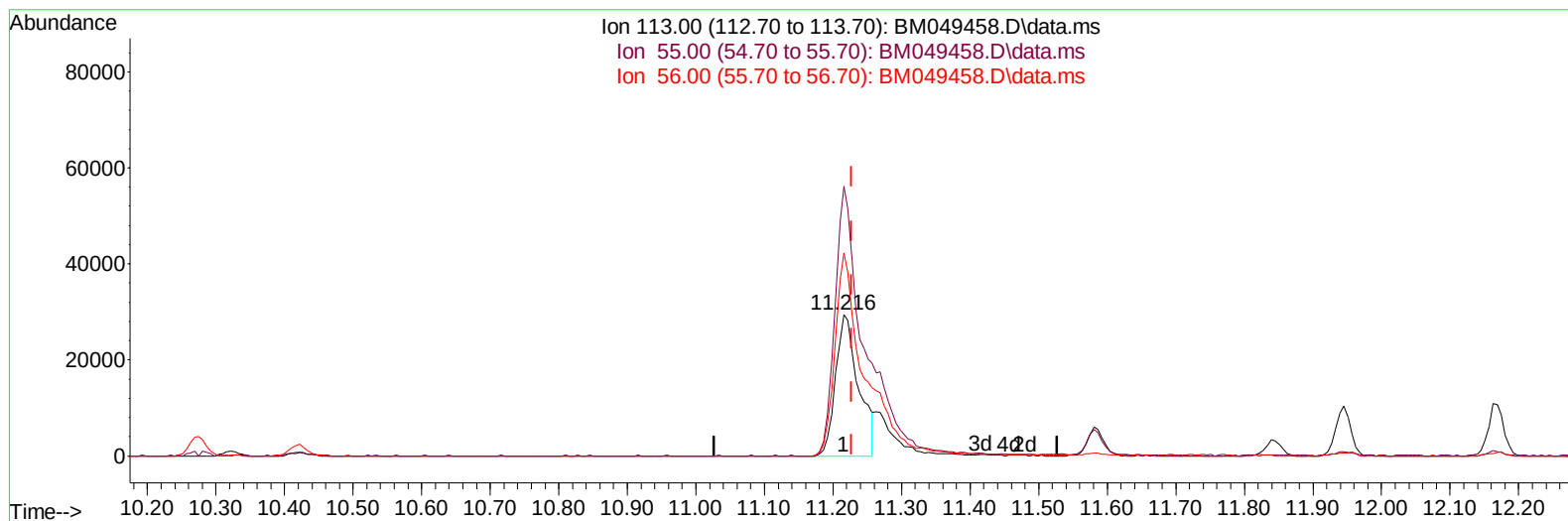
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM012725\  
Data File : BM049458.D  
Acq On : 27 Jan 2025 21:18  
Operator : RC/JU  
Sample : PB166251BS  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SLCS251

Manual IntegrationsAPPROVED

Quant Time: Jan 28 00:47:58 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM011325.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jan 13 16:56:06 2025  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/28/2025  
Supervised By :mohammad ahmed 01/31/2025



TIC: BM049458.D\data.ms

(34) Caprolactam

11.216min (-0.012) 24.02 ng/ul

response 68712

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 193.90 | 190.49 |
| 56.00  | 143.50 | 143.32 |
| 0.00   | 0.00   | 0.00   |

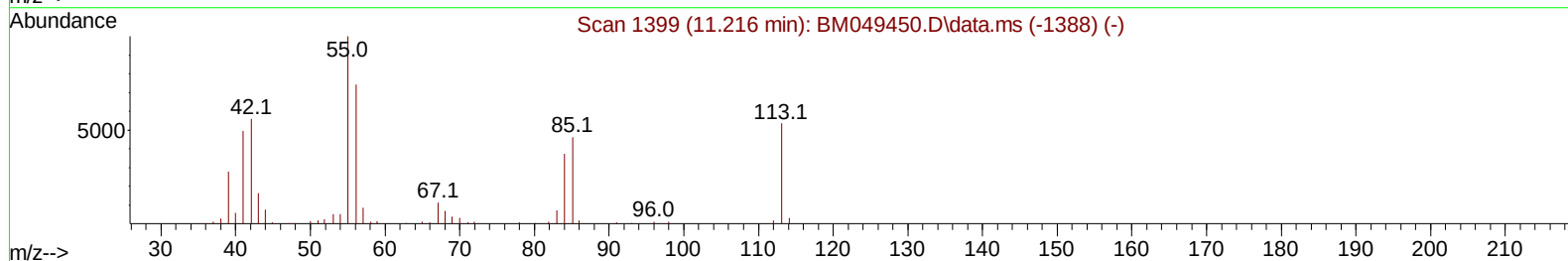
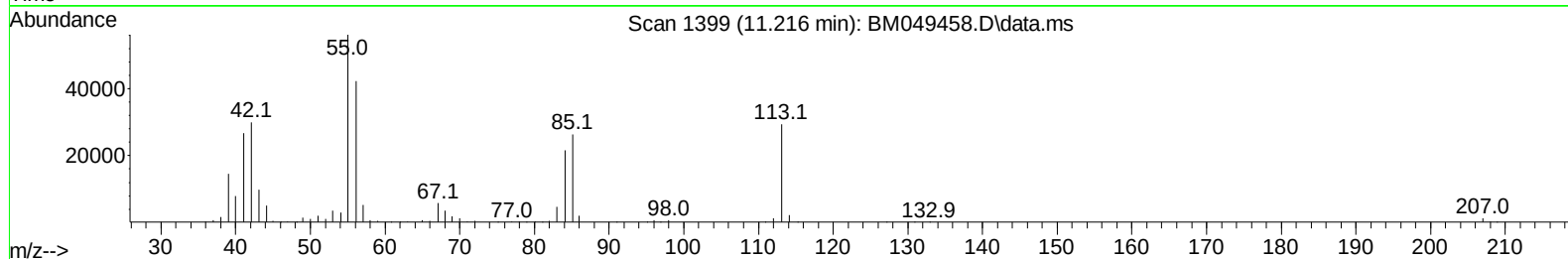
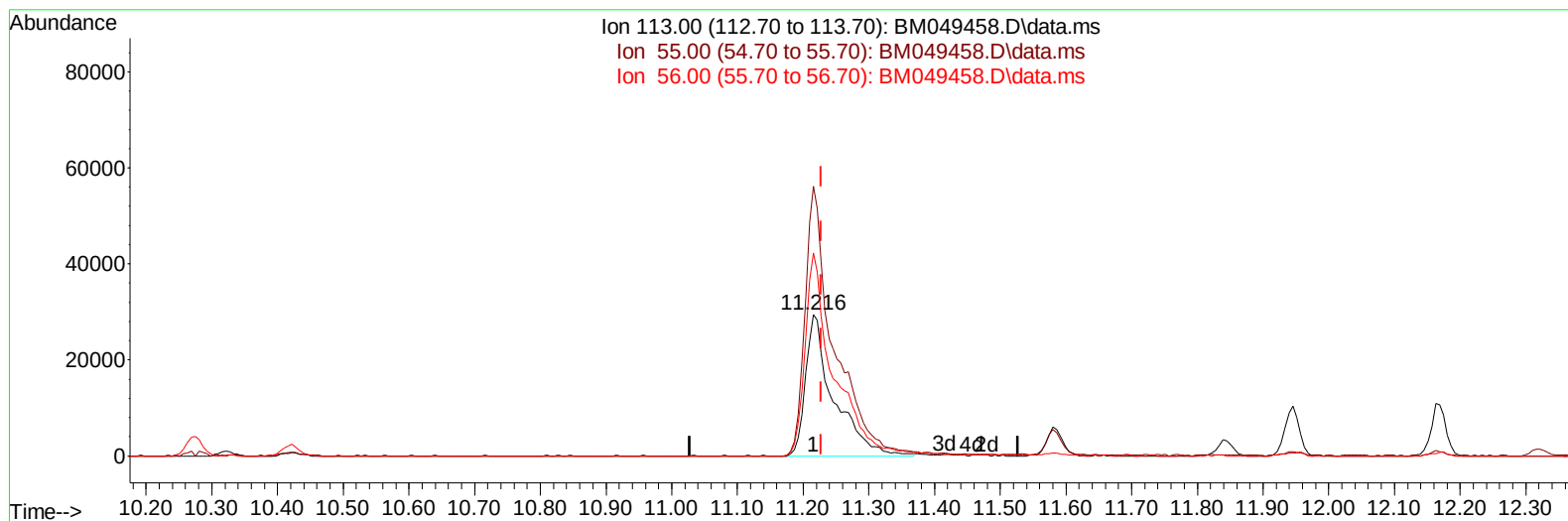
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM012725\  
Data File : BM049458.D  
Acq On : 27 Jan 2025 21:18  
Operator : RC/JU  
Sample : PB166251BS  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SLCS251

Manual IntegrationsAPPROVED

Quant Time: Jan 28 00:49:45 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM011325.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jan 13 16:56:06 2025  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/28/2025  
Supervised By :mohammad ahmed 01/31/2025



TIC: BM049458.D\data.ms

**(34) Caprolactam**

11.216min (-0.012) 30.64 ng/ul m

response 87674

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 193.90 | 190.49 |
| 56.00  | 143.50 | 143.32 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM012725\  
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 Acq On : 27 Jan 2025 21:18  
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 Sample : PB166251BS  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SLCS251

Manual IntegrationsAPPROVED

Quant Time: Jan 28 01:29:55 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM011325.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jan 13 16:56:06 2025  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/28/2025  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.516  | 152  | 144355   | 20.000 | ng/uL  | 0.00     |
| 20) Naphthalene-d8            | 10.275 | 136  | 554528   | 20.000 | ng/uL  | -0.01    |
| 38) Acenaphthene-d10          | 14.157 | 164  | 383480   | 20.000 | ng/uL  | 0.00     |
| 64) Phenanthrene-d10          | 16.904 | 188  | 811079   | 20.000 | ng/uL  | -0.01    |
| 79) Chrysene-d12              | 21.139 | 240  | 774804   | 20.000 | ng/uL  | 0.00     |
| 88) Perylene-d12              | 23.915 | 264  | 758984   | 20.000 | ng/uL  | -0.02    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.122  | 96   | 25056    | 7.182  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.516  | 84   | 373035   | 33.586 | ng/uL  | 0.00     |
| 7) Phenol-d5                  | 6.710  | 99   | 425156   | 35.085 | ng/uL  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.863  | 67   | 262896   | 32.581 | ng/uL  | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.057  | 132  | 328137   | 35.398 | ng/uL  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.228  | 113  | 314750   | 33.495 | ng/uL  | -0.01    |
| 21) Nitrobenzene-d5           | 8.657  | 128  | 139631   | 33.020 | ng/uL  | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.369  | 143  | 156454   | 33.096 | ng/uL  | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 9.910  | 165  | 348062   | 36.083 | ng/uL  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.422 | 131  | 361725   | 28.498 | ng/uL  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.574 | 166  | 955961   | 33.493 | ng/uL  | -0.01    |
| 49) Acenaphthylene-d8         | 13.839 | 160  | 1095652  | 33.626 | ng/uL  | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.380 | 143  | 150555   | 31.854 | ng/uL  | -0.01    |
| 60) Fluorene-d10              | 15.157 | 176  | 848847   | 34.427 | ng/uL  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.286 | 200  | 156351   | 30.247 | ng/uL  | 0.00     |
| 73) Anthracene-d10            | 17.004 | 188  | 1338937  | 33.394 | ng/uL  | -0.01    |
| 81) Pyrene-d10                | 19.309 | 212  | 1522684  | 32.292 | ng/uL  | -0.01    |
| 92) Benzo(a)pyrene-d12        | 23.733 | 264  | 1378758  | 33.840 | ng/uL  | -0.01    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.152  | 88   | 49598    | 12.976 | ng/uL  | 96       |
| 5) Pyridine                   | 3.534  | 79   | 381199   | 33.952 | ng/uL  | 98       |
| 6) Benzaldehyde               | 6.675  | 77   | 22441    | 3.435  | ng/uL  | 93       |
| 8) Phenol                     | 6.740  | 94   | 446334   | 35.426 | ng/uL  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 6.957  | 93   | 325057   | 31.683 | ng/uL  | 99       |
| 12) 2-Chlorophenol            | 7.087  | 128  | 339001   | 35.182 | ng/uL  | 96       |
| 13) 2-Methylphenol            | 7.963  | 108  | 307073   | 33.359 | ng/uL  | 97       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.034  | 45   | 521572   | 33.920 | ng/uL  | 99       |
| 16) Acetophenone              | 8.328  | 105  | 480611   | 31.316 | ng/uL  | 95       |
| 17) N-Nitroso-di-n-propyla... | 8.322  | 70   | 271691   | 32.775 | ng/uL  | 96       |
| 18) 4-Methylphenol            | 8.292  | 108  | 336409   | 33.357 | ng/uL  | 98       |
| 19) Hexachloroethane          | 8.575  | 117  | 131604   | 31.787 | ng/uL  | 98       |
| 22) Nitrobenzene              | 8.698  | 77   | 366200   | 33.427 | ng/uL  | 100      |
| 23) Isophorone                | 9.228  | 82   | 700289   | 32.899 | ng/uL  | 100      |
| 25) 2-Nitrophenol             | 9.404  | 139  | 169467   | 32.618 | ng/uL  | 98       |
| 26) 2,4-Dimethylphenol        | 9.475  | 107  | 331824   | 34.126 | ng/uL  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 9.710  | 93   | 426297   | 32.584 | ng/uL  | 98       |
| 29) 2,4-Dichlorophenol        | 9.933  | 162  | 343637   | 35.756 | ng/uL  | 98       |
| 30) Naphthalene               | 10.322 | 128  | 947157   | 32.184 | ng/uL  | 99       |
| 32) 4-Chloroaniline           | 10.445 | 127  | 192108   | 15.871 | ng/uL  | 99       |
| 33) Hexachlorobutadiene       | 10.616 | 225  | 225009   | 32.791 | ng/uL  | 98       |
| 34) Caprolactam               | 11.216 | 113  | 87674m   | 30.644 | ng/uL  |          |
| 35) 4-Chloro-3-methylphenol   | 11.580 | 107  | 313778   | 34.919 | ng/uL  | 99       |

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## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/28/2025  
 Supervised By :mohammad ahmed 01/31/2025

Quant Time: Jan 28 01:29:55 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM011325.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jan 13 16:56:06 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 11.945 | 142  | 673606   | 31.987 | ng/ul | 100      |
| 37) 1-Methylnaphthalene       | 12.169 | 142  | 679217   | 32.416 | ng/ul | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.322 | 216  | 430380   | 32.423 | ng/ul | 99       |
| 40) Hexachlorocyclopentadiene | 12.304 | 237  | 200127   | 25.129 | ng/ul | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.574 | 196  | 283909   | 34.769 | ng/ul | 97       |
| 42) 2,4,5-Trichlorophenol     | 12.645 | 196  | 307451   | 35.130 | ng/ul | 99       |
| 43) 1,1'-Biphenyl             | 12.980 | 154  | 912504   | 32.500 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.016 | 162  | 741023   | 33.002 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.233 | 65   | 210494   | 35.528 | ng/ul | 99       |
| 47) Dimethylphthalate         | 13.621 | 163  | 942529   | 33.382 | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene        | 13.739 | 165  | 182341   | 34.616 | ng/ul | 99       |
| 50) Acenaphthylene            | 13.869 | 152  | 1111027  | 33.260 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.069 | 138  | 161650   | 31.278 | ng/ul | 95       |
| 52) Acenaphthene              | 14.216 | 153  | 788230   | 33.027 | ng/ul | 99       |
| 53) 2,4-Dinitrophenol         | 14.280 | 184  | 91900    | 27.857 | ng/ul | 96       |
| 55) 4-Nitrophenol             | 14.398 | 109  | 112773   | 32.356 | ng/ul | 99       |
| 56) Dibenzofuran              | 14.557 | 168  | 1149918  | 33.929 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 14.533 | 165  | 272507   | 34.862 | ng/ul | 100      |
| 58) 2,3,4,6-Tetrachlorophenol | 14.792 | 232  | 253531   | 34.293 | ng/ul | 98       |
| 59) Diethylphthalate          | 15.004 | 149  | 924665   | 33.604 | ng/ul | 99       |
| 61) Fluorene                  | 15.210 | 166  | 968884   | 34.989 | ng/ul | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.215 | 204  | 513708   | 34.040 | ng/ul | 97       |
| 63) 4-Nitroaniline            | 15.239 | 138  | 187588   | 40.341 | ng/ul | 99       |
| 66) 4,6-Dinitro-2-methylph... | 15.298 | 198  | 173369   | 30.760 | ng/ul | 95       |
| 67) N-Nitrosodiphenylamine    | 15.427 | 169  | 782407   | 32.512 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.110 | 248  | 302023   | 32.748 | ng/ul | 96       |
| 69) Hexachlorobenzene         | 16.215 | 284  | 347767   | 32.422 | ng/ul | 98       |
| 70) Atrazine                  | 16.392 | 200  | 285847   | 29.053 | ng/ul | 98       |
| 71) Pentachlorophenol         | 16.562 | 266  | 207958   | 29.265 | ng/ul | 99       |
| 72) Phenanthrene              | 16.951 | 178  | 1490799  | 33.062 | ng/ul | 99       |
| 74) Anthracene                | 17.039 | 178  | 1511894  | 33.405 | ng/ul | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 12.939 | 216  | 424475   | 31.314 | ng/uL | 98       |
| 76) Pentachlorobenzene        | 14.474 | 250  | 413200   | 30.539 | ng/uL | 100      |
| 77) Carbazole                 | 17.315 | 167  | 1324413  | 32.982 | ng/ul | 100      |
| 78) Di-n-butylphthalate       | 17.898 | 149  | 1589108  | 32.298 | ng/ul | 100      |
| 80) Fluoranthene              | 18.974 | 202  | 1747984  | 32.260 | ng/ul | 99       |
| 82) Pyrene                    | 19.339 | 202  | 1867719  | 32.382 | ng/ul | 98       |
| 83) Butylbenzylphthalate      | 20.268 | 149  | 640951   | 31.389 | ng/ul | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.062 | 252  | 454612   | 31.333 | ng/ul | 99       |
| 85) Benzo(a)anthracene        | 21.121 | 228  | 1786921  | 33.094 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.074 | 149  | 965114   | 31.780 | ng/ul | 100      |
| 87) Chrysene                  | 21.180 | 228  | 1665191  | 33.781 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 22.139 | 149  | 1514231  | 27.386 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.044 | 252  | 1642088  | 32.324 | ng/ul | 100      |
| 91) Benzo(k)fluoranthene      | 23.103 | 252  | 1663466  | 32.841 | ng/ul | 100      |
| 93) Benzo(a)pyrene            | 23.791 | 252  | 1548264  | 33.741 | ng/ul | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 26.997 | 276  | 1964830  | 37.328 | ng/ul | 99       |
| 95) Dibenzo(a,h)anthracene    | 27.056 | 278  | 1600343  | 37.328 | ng/ul | 99       |
| 96) Benzo(g,h,i)perylene      | 27.962 | 276  | 1496335  | 38.081 | ng/ul | 100      |

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

**Instrument :**

BNA\_M

**ClientSampleId :**

SLCS251

**Manual IntegrationsAPPROVED**

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