

(QT Reviewed)

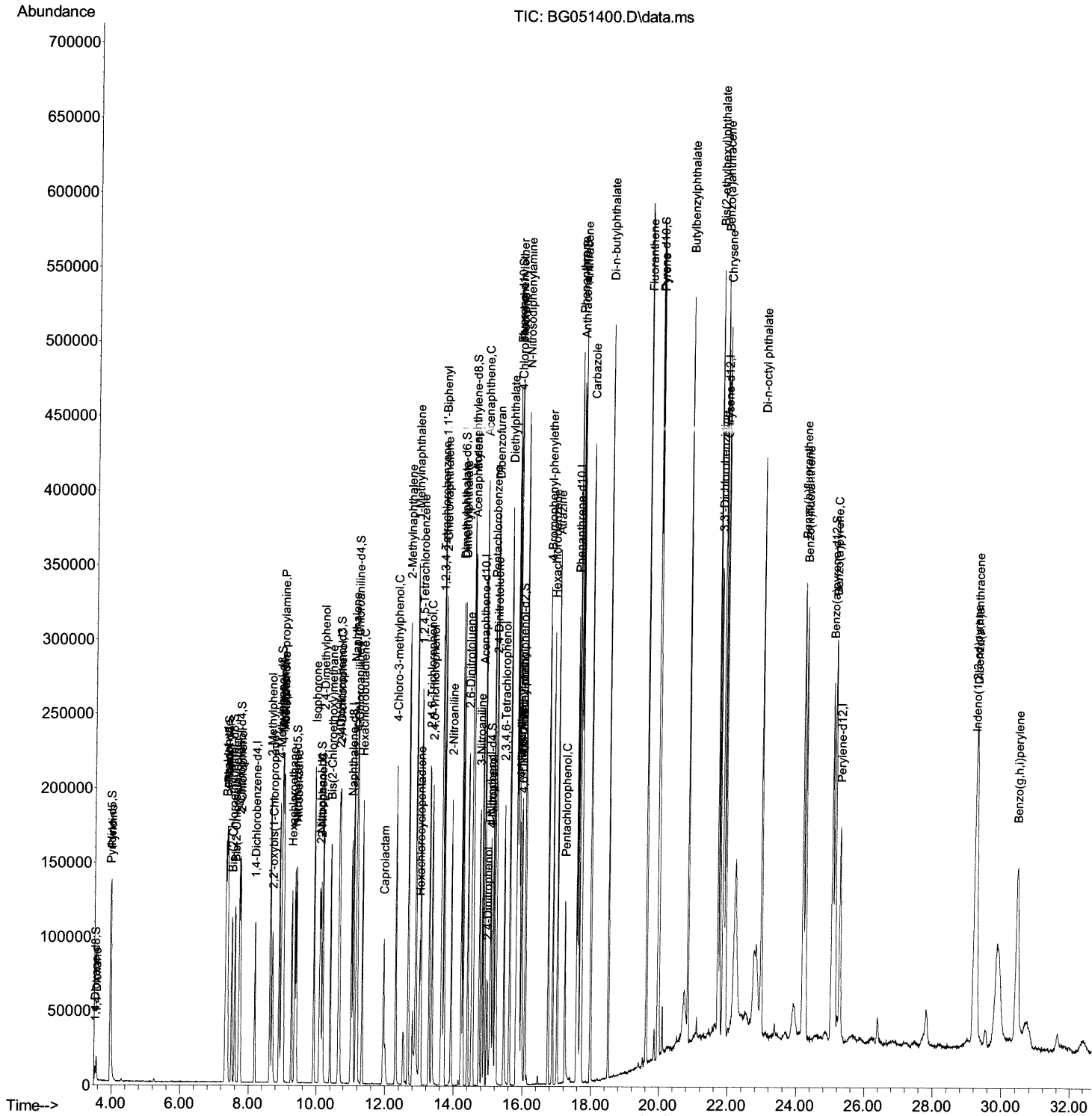
```
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051400.D
Acq On    : 8 Dec 2021    2:49
Operator  : CG/JU
Sample    : PB141220BS
Misc      :
ALS Vial  : 58    Sample Multiplier: 1
```

Instrument :
BNA_G
ClientSampleId :
SLCS220

Manual IntegrationsAPPROVED

Quant Time: Dec 08 05:31:17 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 03 15:23:09 2021
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/08/2021
Supervised By :mohammad ahmed 12/15/2021



Quantitation Report (Qedit)

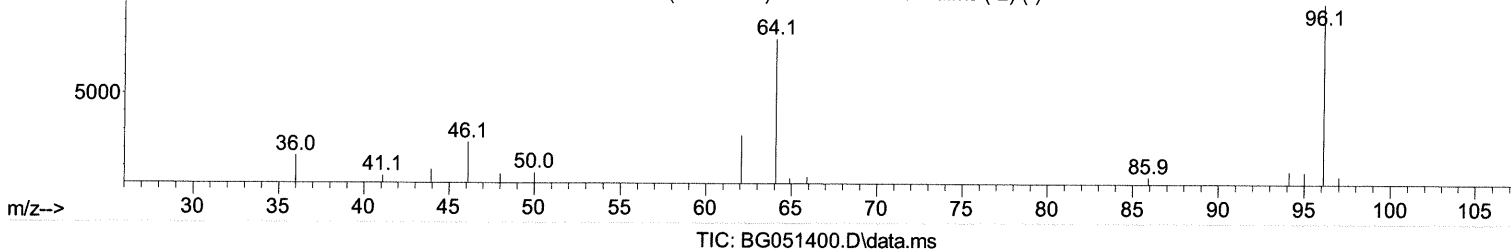
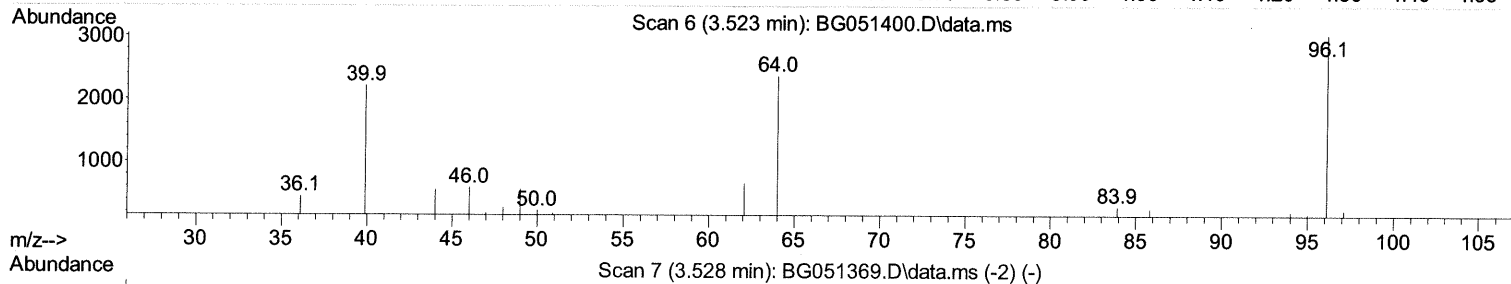
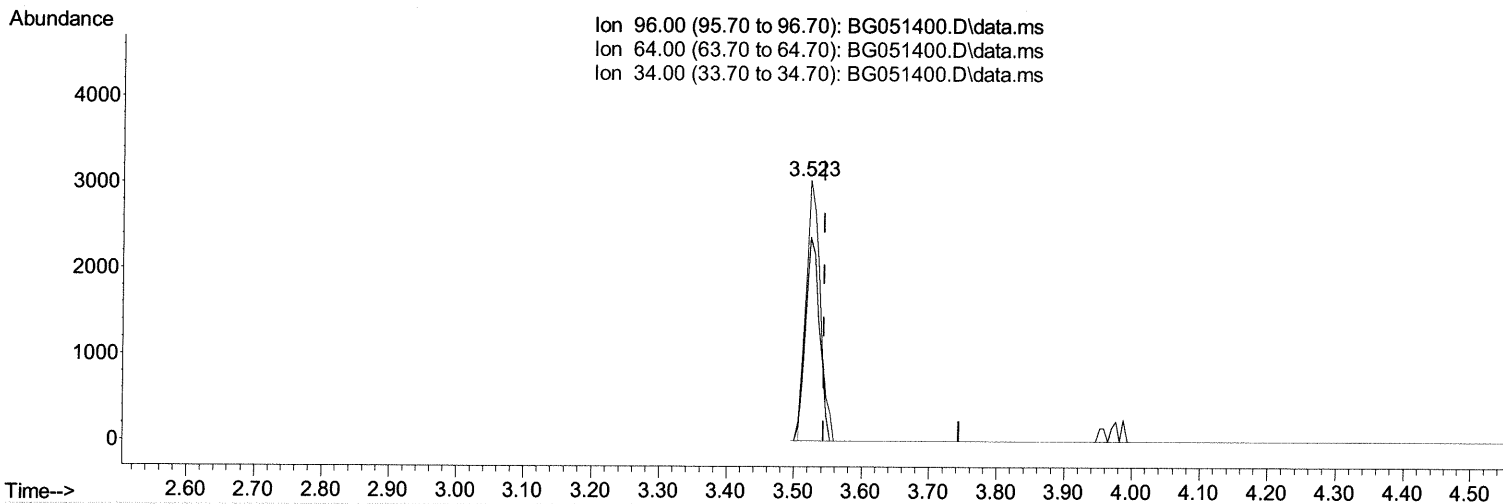
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(3) 1,4-Dioxane-d8 (S)

3.523min (-0.021) 5.05 ng/uL

response 4395

| Ion | Exp% | Act% |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 77.60 | 78.13 |
| 34.00 | 0.00 | 0.00 |
| 0.00 | 0.00 | 0.00 |

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 Operator : CG/JU
 Sample : PB141220BS
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Instrument :

BNA_G

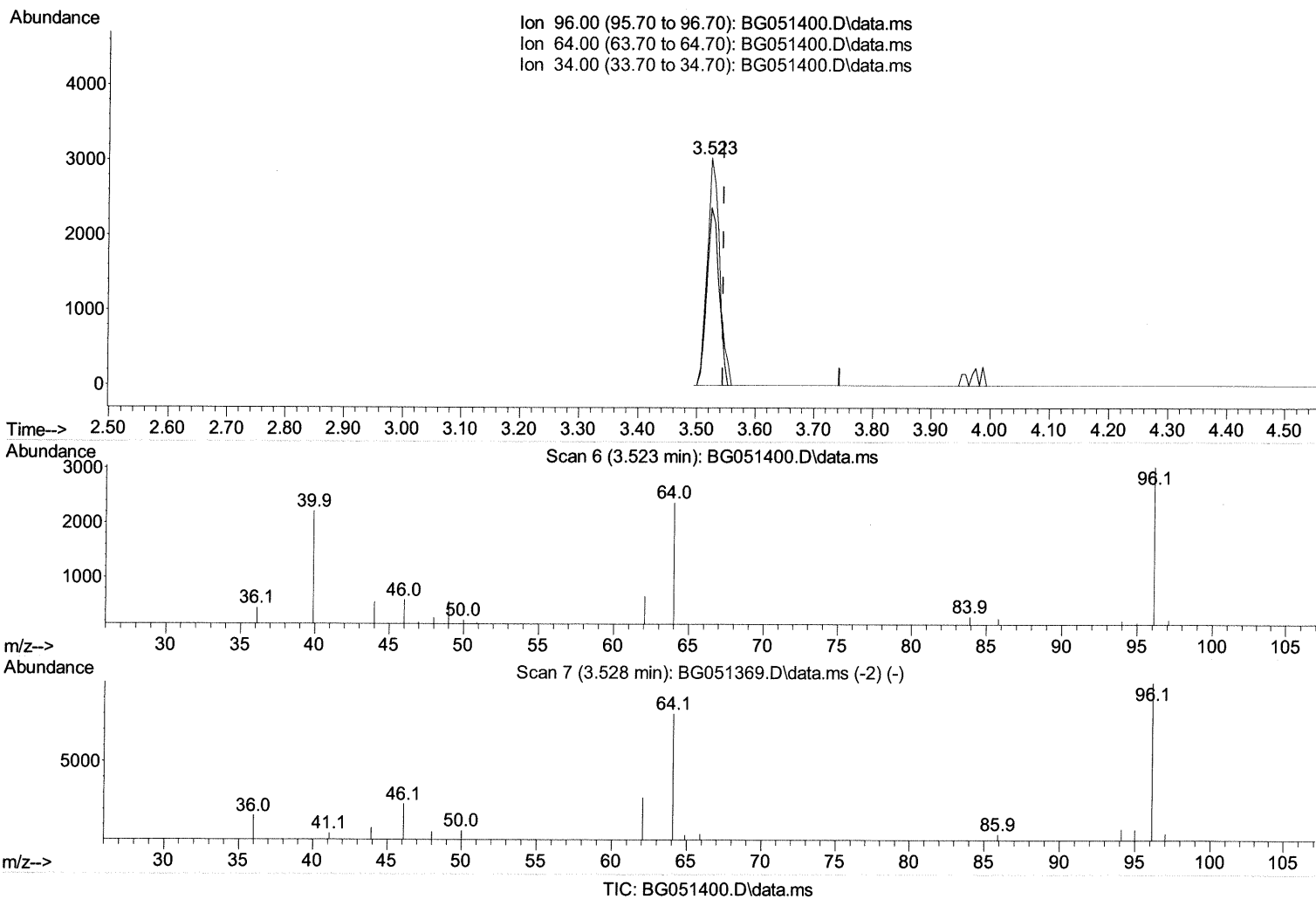
ClientSampleId :

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(3) 1,4-Dioxane-d8 (S)

3.523min (-0.021) 5.14 ng/uL m 12/16/21 JU

response 4478

| Ion | Exp% | Act% |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 77.60 | 78.13 |
| 34.00 | 0.00 | 0.00 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

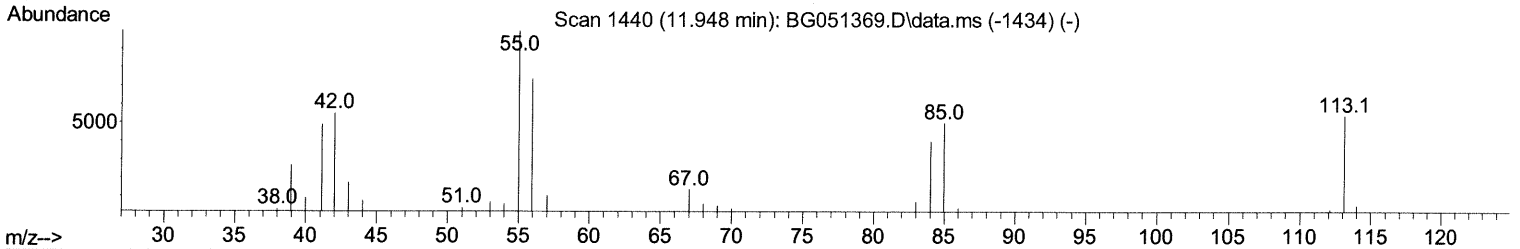
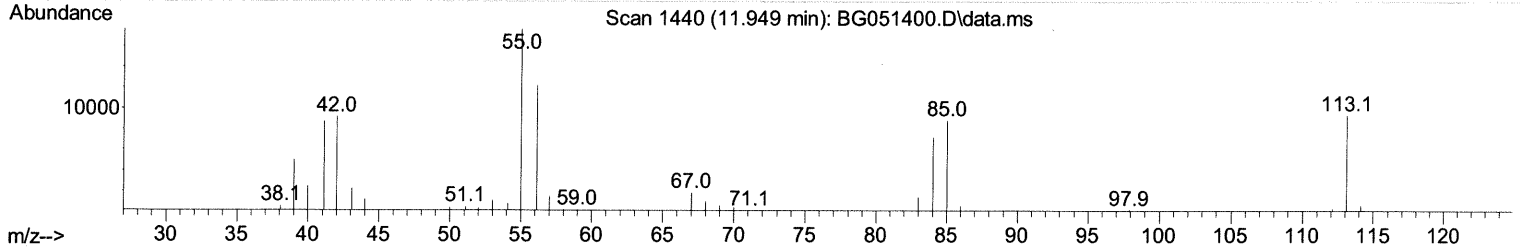
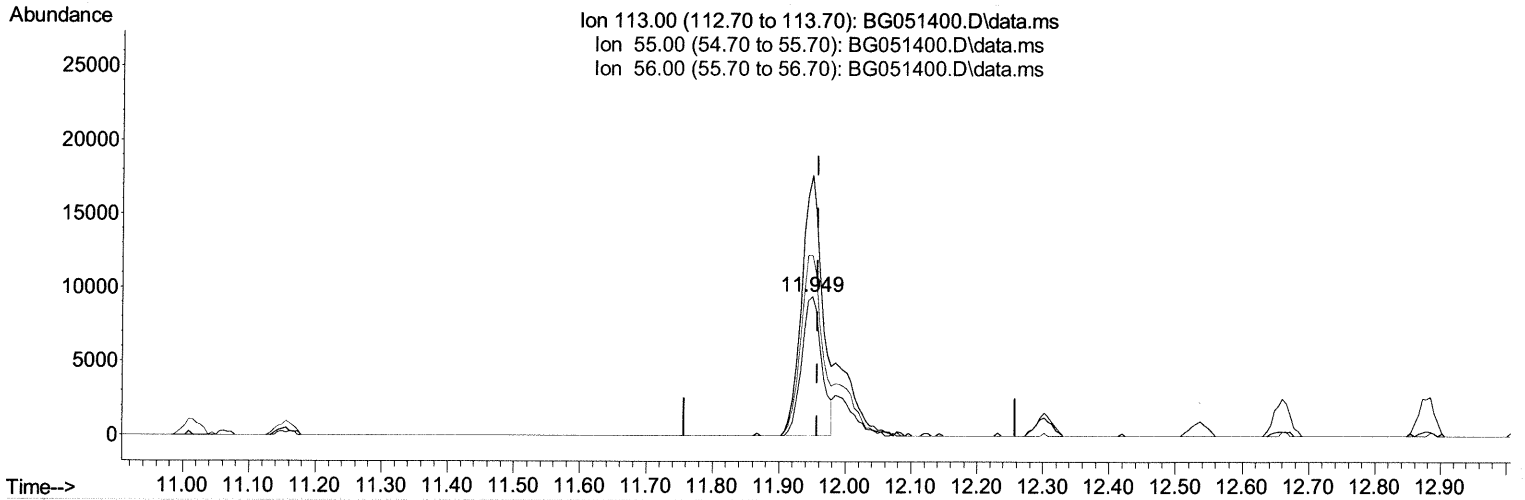
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 Operator : CG/JU
 Sample : PB141220BS
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
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TIC: BG051400.D\data.ms

(34) Caprolactam

11.949min (-0.009) 23.46 ng/ul

response 20091

| Ion | Exp% | Act% |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 183.80 | 187.20 |
| 56.00 | 136.50 | 129.74 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

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 Operator : CG/JU
 Sample : PB141220BS
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :

BNA_G

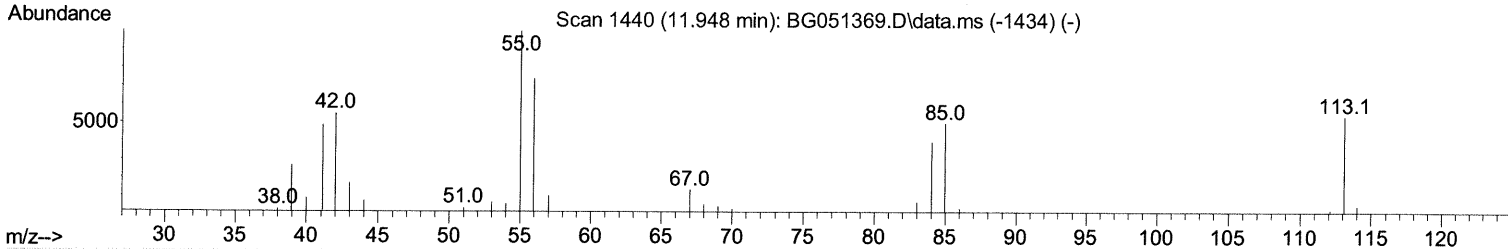
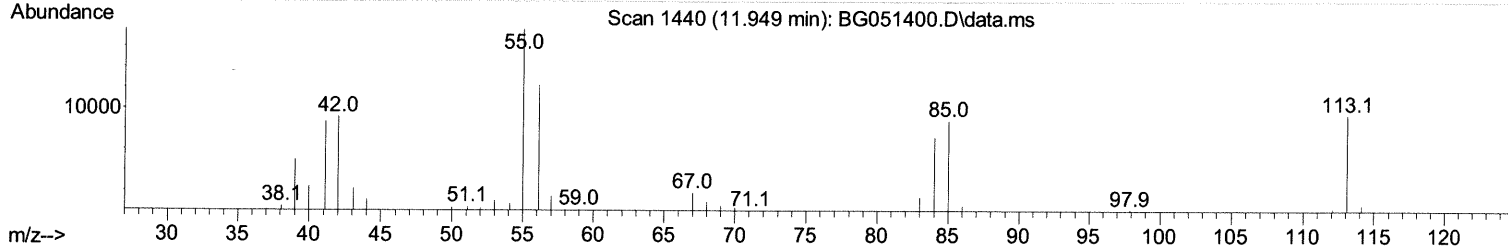
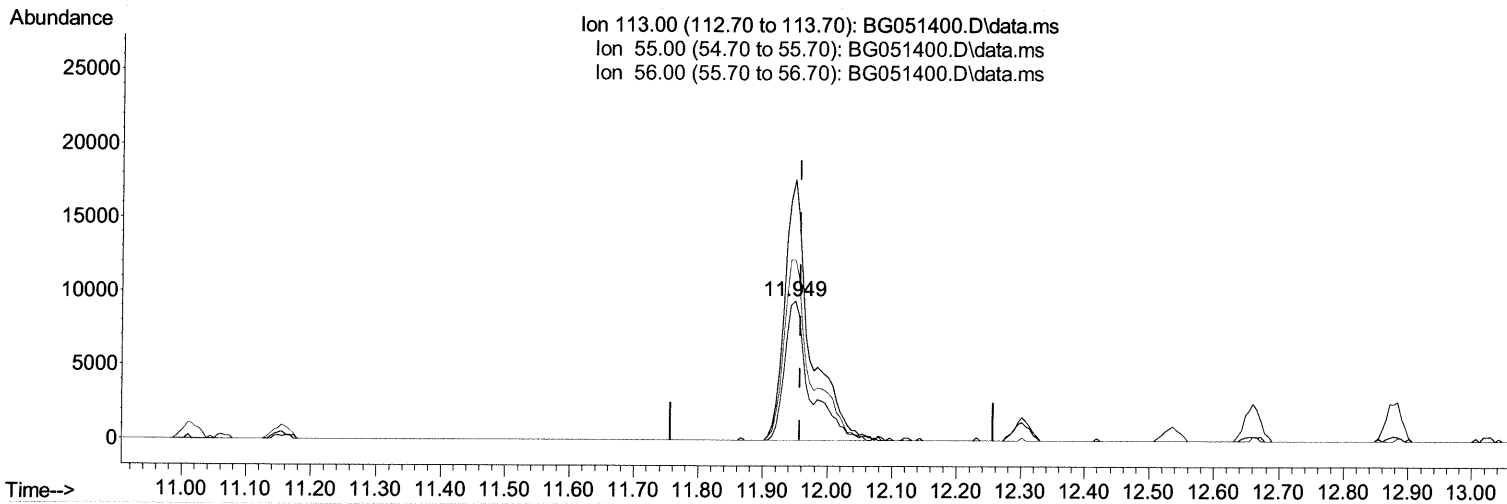
ClientSampleId :

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TIC: BG051400.D\data.ms

(34) Caprolactam

11.949min (-0.009) 30.26 ng/ul m 12/11/2021

response 25917

| Ion | Exp% | Act% |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 183.80 | 187.20 |
| 56.00 | 136.50 | 129.74 |
| 0.00 | 0.00 | 0.00 |

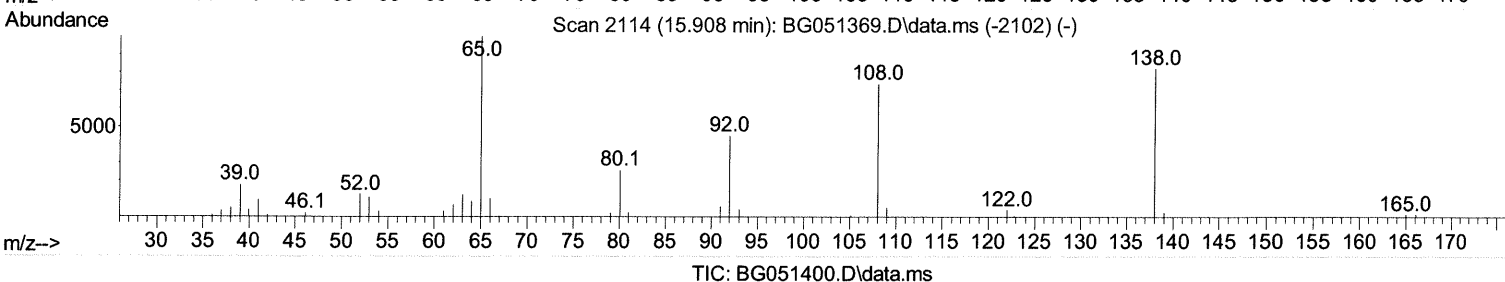
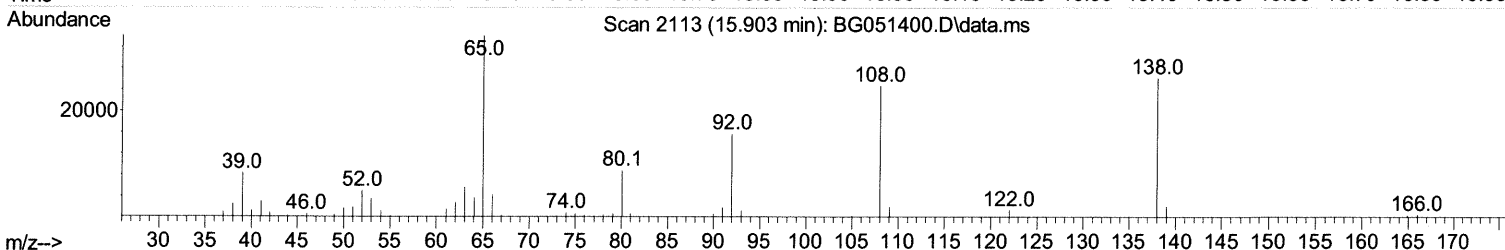
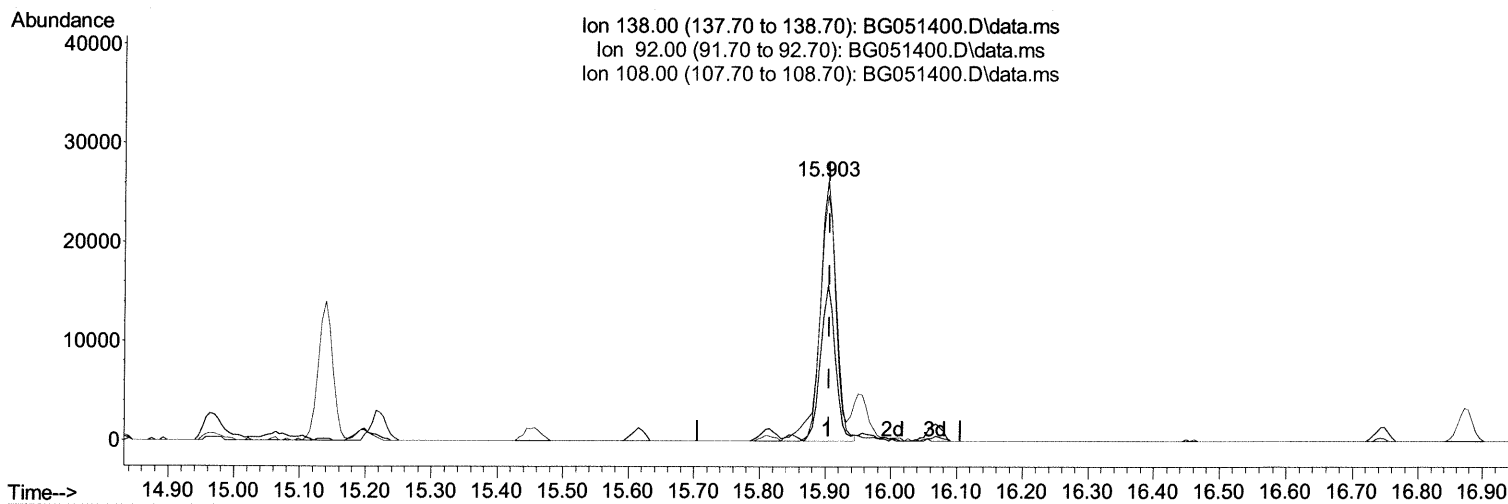
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051400.D
Acq On : 8 Dec 2021 2:49
Operator : CG/JU
Sample : PB141220BS
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS220

Manual IntegrationsAPPROVED

Quant Time: Dec 08 05:31:17 2021
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Reviewed By :Jagrut Upadhyay 12/08/2021
Supervised By :mohammad ahmed 12/15/2021



(63) 4-Nitroaniline

15.903min (-0.003) 32.66 ng/ul

response 45945

| Ion | Exp% | Act% |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00 | 61.60 | 59.84 |
| 108.00 | 90.70 | 94.34 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

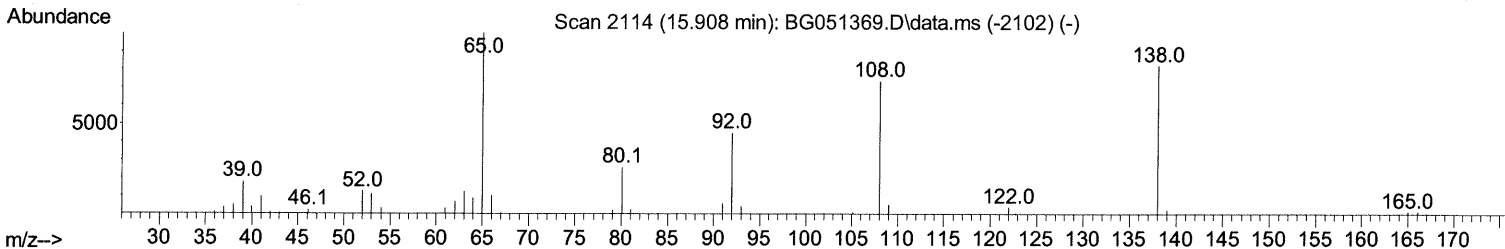
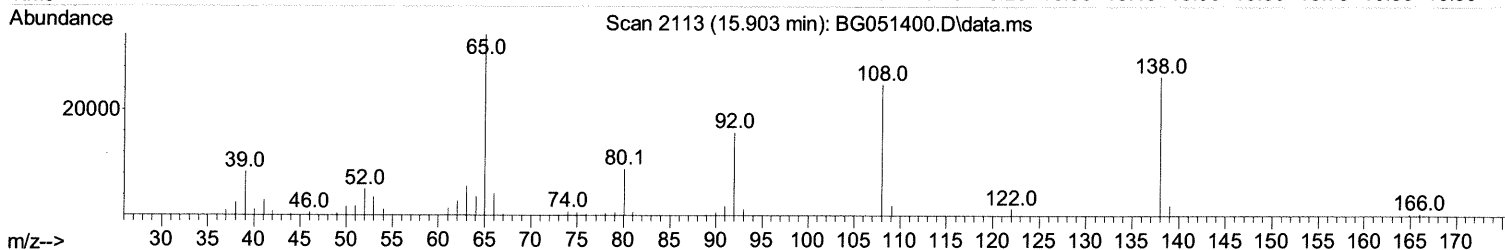
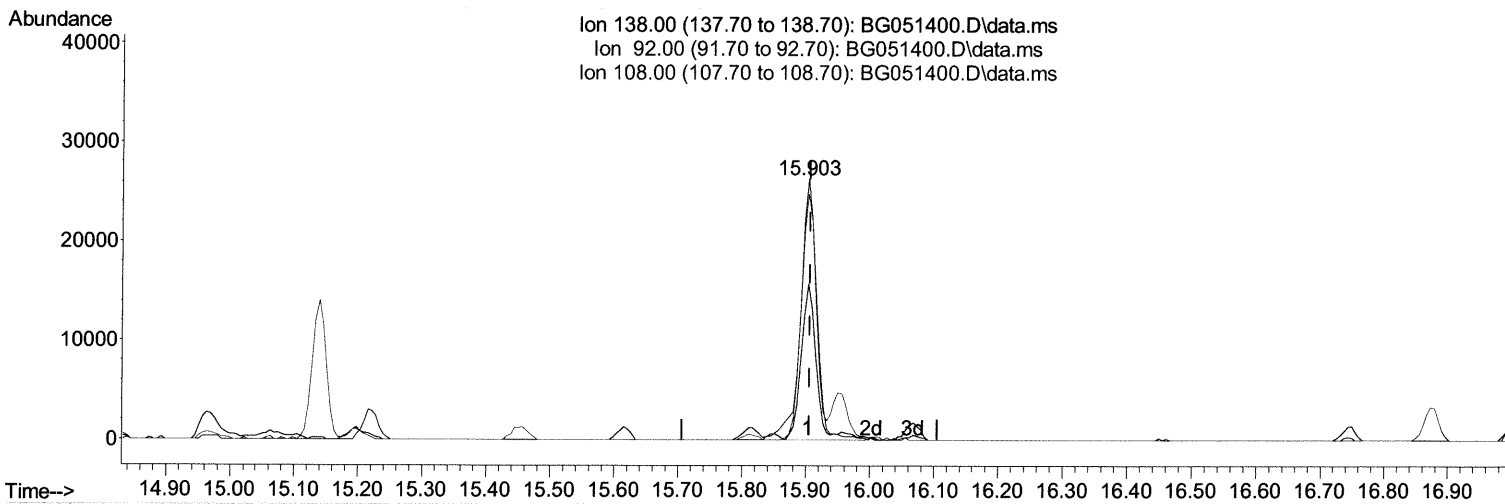
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
 Data File : BG051400.D
 Acq On : 8 Dec 2021 2:49
 Operator : CG/JU
 Sample : PB141220B5
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLCS220

Manual IntegrationsAPPROVED

Quant Time: Dec 08 05:31:17 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 03 15:23:09 2021
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/08/2021
 Supervised By :mohammad ahmed 12/15/2021



TIC: BG051400.D\data.ms

(63) 4-Nitroaniline

15.903min (-0.003) 33.26 ng/ul m (21/6/2130)

response 46779

| Ion | Exp% | Act% |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00 | 61.60 | 59.84 |
| 108.00 | 90.70 | 94.34 |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
 Data File : BG051400.D
 Acq On : 8 Dec 2021 2:49
 Operator : CG/JU
 Sample : PB141220BS
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLCS220

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021
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| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|-------------------|
| ----- | | | | | | |
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 8.189 | 152 | 30263 | 20.000 | ng/ul | -0.01 |
| 20) Naphthalene-d8 | 11.015 | 136 | 136990 | 20.000 | ng/ul | -0.01 |
| 38) Acenaphthene-d10 | 14.822 | 164 | 89384 | 20.000 | ng/ul | 0.00 |
| 64) Phenanthrene-d10 | 17.572 | 188 | 190863 | 20.000 | ng/ul | 0.00 |
| 79) Chrysene-d12 | 21.873 | 240 | 152888 | 20.000 | ng/ul | 0.00 |
| 88) Perylene-d12 | 25.274 | 264 | 153814 | 20.000 | ng/ul | 0.00 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.523 | 96 | 4478m | 5.142 | ng/ul | -0.02 12/16/21 JU |
| 4) Pyridine-d5 | 3.952 | 84 | 64369 | 25.189 | ng/ul | -0.03 |
| 7) Phenol-d5 | 7.354 | 99 | 92084 | 30.787 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.501 | 67 | 57045 | 30.367 | ng/ul | -0.01 |
| 11) 2-Chlorophenol-d4 | 7.719 | 132 | 66656 | 30.948 | ng/ul | -0.01 |
| 15) 4-Methylphenol-d8 | 8.905 | 113 | 73909 | 30.621 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5 | 9.370 | 128 | 35120 | 30.370 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4 | 10.092 | 143 | 41243 | 31.617 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.645 | 165 | 70254 | 31.742 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4 | 11.156 | 131 | 128289 | 39.614 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6 | 14.217 | 166 | 209508 | 30.462 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8 | 14.522 | 160 | 271495 | 31.305 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4 | 15.051 | 143 | 28833 | 25.900 | ng/ul | 0.00 |
| 60) Fluorene-d10 | 15.815 | 176 | 190544 | 30.766 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.956 | 200 | 30014 | 25.484 | ng/ul | 0.00 |
| 73) Anthracene-d10 | 17.672 | 188 | 283450 | 31.052 | ng/ul | 0.00 |
| 81) Pyrene-d10 | 19.951 | 212 | 301821 | 32.626 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12 | 25.039 | 264 | 262370 | 31.939 | ng/ul | 0.00 |
| Target Compounds | | | | | | |
| | | | | | Qvalue | |
| 2) 1,4-Dioxane | 3.565 | 88 | 8983 | 9.146 | ng/ul# | 94 |
| 5) Pyridine | 3.976 | 79 | 66801 | 25.122 | ng/ul | 98 |
| 6) Benzaldehyde | 7.325 | 77 | 62206 | 32.658 | ng/ul | 95 |
| 8) Phenol | 7.378 | 94 | 91857 | 29.646 | ng/ul | 99 |
| 10) Bis(2-Chloroethyl)ether | 7.595 | 93 | 68008 | 29.012 | ng/ul | 99 |
| 12) 2-Chlorophenol | 7.754 | 128 | 65478 | 29.833 | ng/ul | 97 |
| 13) 2-Methylphenol | 8.641 | 108 | 67872 | 29.408 | ng/ul | 95 |
| 14) 2,2'-oxybis(1-Chloropr... | 8.711 | 45 | 102319 | 30.248 | ng/ul | 99 |
| 16) Acetophenone | 9.017 | 105 | 107625 | 28.828 | ng/ul | 98 |
| 17) N-Nitroso-di-n-propyla... | 8.993 | 70 | 62770 | 29.258 | ng/ul | 99 |
| 18) 4-Methylphenol | 8.970 | 108 | 72799 | 29.498 | ng/ul | 94 |
| 19) Hexachloroethane | 9.270 | 117 | 27242 | 29.386 | ng/ul | 96 |
| 22) Nitrobenzene | 9.411 | 77 | 92921 | 30.645 | ng/ul | 96 |
| 23) Isophorone | 9.928 | 82 | 172035 | 29.203 | ng/ul | 99 |
| 25) 2-Nitrophenol | 10.122 | 139 | 40572 | 30.028 | ng/ul | 97 |
| 26) 2,4-Dimethylphenol | 10.174 | 107 | 85071 | 30.795 | ng/ul | 95 |
| 27) Bis(2-Chloroethoxy)met... | 10.404 | 93 | 97353 | 29.935 | ng/ul | 98 |
| 29) 2,4-Dichlorophenol | 10.668 | 162 | 64960 | 29.816 | ng/ul | 96 |
| 30) Naphthalene | 11.068 | 128 | 215012 | 28.846 | ng/ul | 99 |
| 32) 4-Chloroaniline | 11.179 | 127 | 86179 | 26.507 | ng/ul | 98 |
| 33) Hexachlorobutadiene | 11.326 | 225 | 42499 | 28.281 | ng/ul | 98 |
| 34) Caprolactam | 11.949 | 113 | 25917m | 30.259 | ng/ul | > 12/16/21 JU |
| 35) 4-Chloro-3-methylphenol | 12.301 | 107 | 79219 | 30.269 | ng/ul | 97 |

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|-------------------------------|--------|------|----------|--------|--------|--------------|
| 36) 2-Methylnaphthalene | 12.660 | 142 | 147750 | 29.142 | ng/ul | 99 |
| 37) 1-Methylnaphthalene | 12.877 | 142 | 151942 | 29.129 | ng/ul | 97 |
| 39) 1,2,4,5-Tetrachloroben... | 13.024 | 216 | 84950 | 30.273 | ng/ul | 97 |
| 40) Hexachlorocyclopentadiene | 12.989 | 237 | 25774 | 22.724 | ng/ul | 99 |
| 41) 2,4,6-Trichlorophenol | 13.271 | 196 | 54368 | 30.874 | ng/ul | 93 |
| 42) 2,4,5-Trichlorophenol | 13.353 | 196 | 58075 | 31.493 | ng/ul | 99 |
| 43) 1,1'-Biphenyl | 13.653 | 154 | 200383 | 30.015 | ng/ul | 99 |
| 44) 2-Chloronaphthalene | 13.706 | 162 | 156302 | 29.432 | ng/ul | 97 |
| 45) 2-Nitroaniline | 13.917 | 65 | 60070 | 32.683 | ng/ul | 94 |
| 47) Dimethylphthalate | 14.264 | 163 | 203421 | 29.221 | ng/ul | 99 |
| 48) 2,6-Dinitrotoluene | 14.405 | 165 | 44450 | 30.397 | ng/ul | 94 |
| 50) Acenaphthylene | 14.552 | 152 | 256614 | 29.949 | ng/ul | 99 |
| 51) 3-Nitroaniline | 14.740 | 138 | 45677 | 31.601 | ng/ul | 95 |
| 52) Acenaphthene | 14.887 | 153 | 168577 | 29.833 | ng/ul | 97 |
| 53) 2,4-Dinitrophenol | 14.969 | 184 | 20201 | 24.993 | ng/ul# | 82 |
| 55) 4-Nitrophenol | 15.069 | 109 | 24869 | 25.751 | ng/ul | 97 |
| 56) Dibenzofuran | 15.222 | 168 | 238866 | 29.307 | ng/ul | 98 |
| 57) 2,4-Dinitrotoluene | 15.198 | 165 | 63135 | 30.229 | ng/ul | 99 |
| 58) 2,3,4,6-Tetrachlorophenol | 15.451 | 232 | 42622 | 29.433 | ng/ul | 98 |
| 59) Diethylphthalate | 15.615 | 149 | 216294 | 29.600 | ng/ul | 99 |
| 61) Fluorene | 15.868 | 166 | 190740 | 29.216 | ng/ul | 99 |
| 62) 4-Chlorophenyl-phenyle... | 15.850 | 204 | 101433 | 28.829 | ng/ul | 97 |
| 63) 4-Nitroaniline | 15.903 | 138 | 46779m | 33.257 | ng/ul | > 12/16/2024 |
| 66) 4,6-Dinitro-2-methylph... | 15.968 | 198 | 27943 | 24.601 | ng/ul | 95 |
| 67) N-Nitrosodiphenylamine | 16.068 | 169 | 172203 | 31.516 | ng/ul | 99 |
| 68) 4-Bromophenyl-phenylether | 16.749 | 248 | 62143 | 30.379 | ng/ul | 93 |
| 69) Hexachlorobenzene | 16.878 | 284 | 64146 | 30.753 | ng/ul | 96 |
| 70) Atrazine | 17.014 | 200 | 68333 | 29.757 | ng/ul | 98 |
| 71) Pentachlorophenol | 17.231 | 266 | 26953 | 29.162 | ng/ul | 96 |
| 72) Phenanthrene | 17.619 | 178 | 317389 | 30.118 | ng/ul | 99 |
| 74) Anthracene | 17.707 | 178 | 317149 | 30.302 | ng/ul | 98 |
| 75) 1,2,3,4-Tetrachloroben... | 13.629 | 216 | 87837 | 31.551 | ng/ul | 97 |
| 76) Pentachlorobenzene | 15.139 | 250 | 78992 | 30.452 | ng/ul | 99 |
| 77) Carbazole | 17.983 | 167 | 279647 | 30.440 | ng/ul | 99 |
| 78) Di-n-butylphthalate | 18.500 | 149 | 365084 | 30.820 | ng/ul | 99 |
| 80) Fluoranthene | 19.622 | 202 | 367379 | 32.333 | ng/ul# | 95 |
| 82) Pyrene | 19.981 | 202 | 349702 | 31.464 | ng/ul | 97 |
| 83) Butylbenzylphthalate | 20.838 | 149 | 145832 | 31.561 | ng/ul | 97 |
| 84) 3,3'-Dichlorobenzidine | 21.761 | 252 | 110510 | 31.045 | ng/ul | 100 |
| 85) Benzo(a)anthracene | 21.855 | 228 | 319265 | 30.788 | ng/ul | 99 |
| 86) Bis(2-ethylhexyl)phtha... | 21.708 | 149 | 209352 | 31.486 | ng/ul | 100 |
| 87) Chrysene | 21.925 | 228 | 304553 | 30.572 | ng/ul | 99 |
| 89) Di-n-octyl phthalate | 22.971 | 149 | 347736 | 31.206 | ng/ul | 100 |
| 90) Benzo(b)fluoranthene | 24.187 | 252 | 316116 | 30.453 | ng/ul | 99 |
| 91) Benzo(k)fluoranthene | 24.258 | 252 | 291327 | 29.907 | ng/ul | 99 |
| 93) Benzo(a)pyrene | 25.116 | 252 | 304470 | 30.745 | ng/ul | 98 |
| 94) Indeno(1,2,3-cd)pyrene | 29.205 | 276 | 335968 | 30.317 | ng/ul | 99 |
| 95) Dibenzo(a,h)anthracene | 29.246 | 278 | 285145 | 30.330 | ng/ul | 99 |
| 96) Benzo(g,h,i)perylene | 30.433 | 276 | 271108 | 29.077 | ng/ul | 99 |

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