

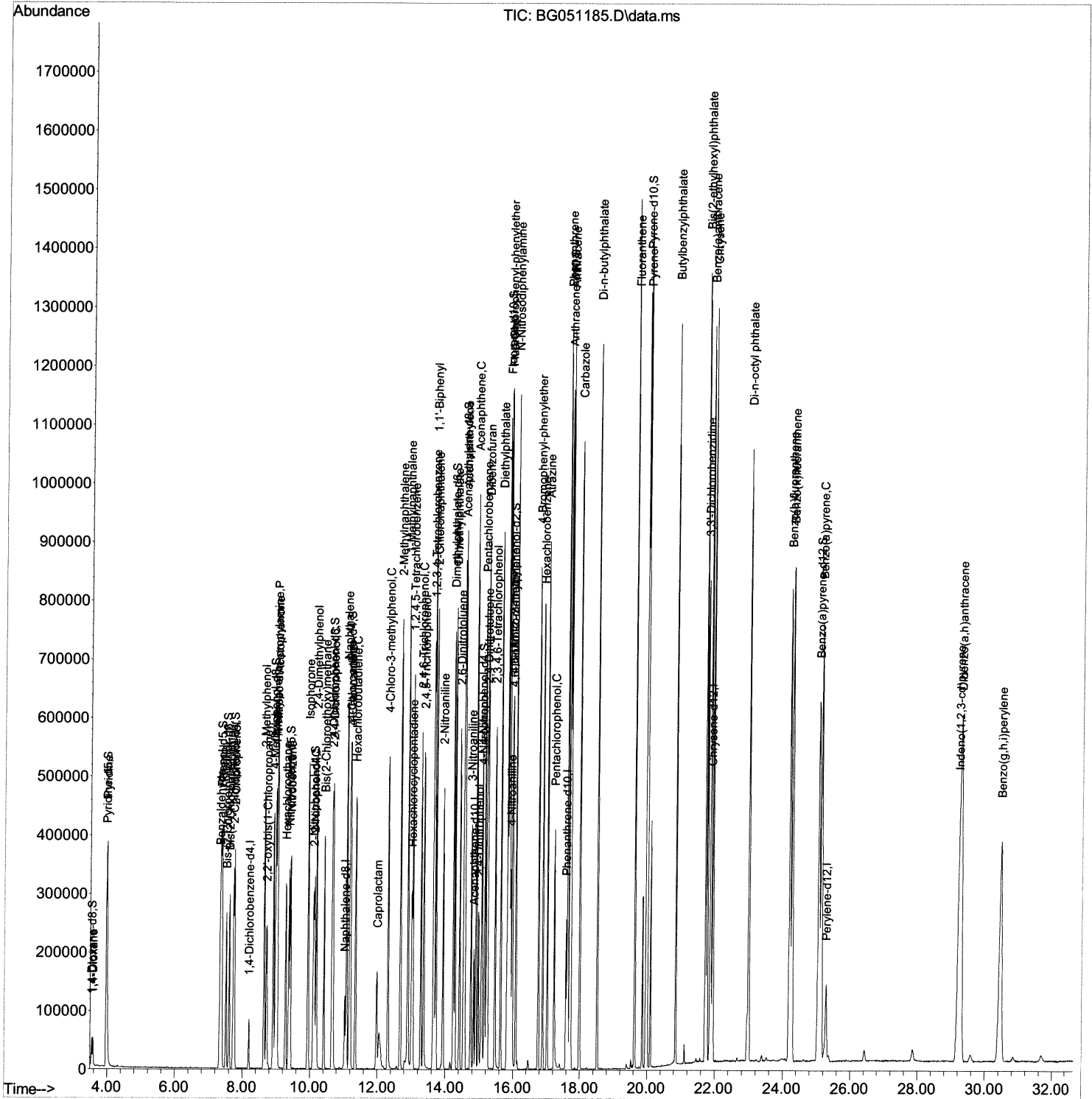
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Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\  
Data File : BG051185.D  
Acq On    : 23 Nov 2021  13:37  
Operator  : CG/JU  
Sample    : SSTD08023  
Misc      :  
ALS Vial  : 6    Sample Multiplier: 1
```

Instrument :
BNA_G
ClientSampleId :
SSTD080424

Manual IntegrationsAPPROVED

Quant Time: Nov 23 14:56:53 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 23 13:54:17 2021
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/23/2021
Supervised By :mohammad ahmed 11/30/2021



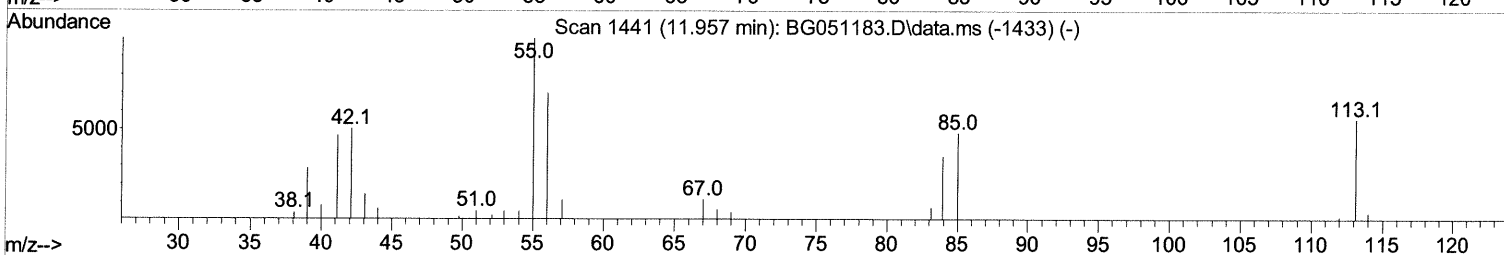
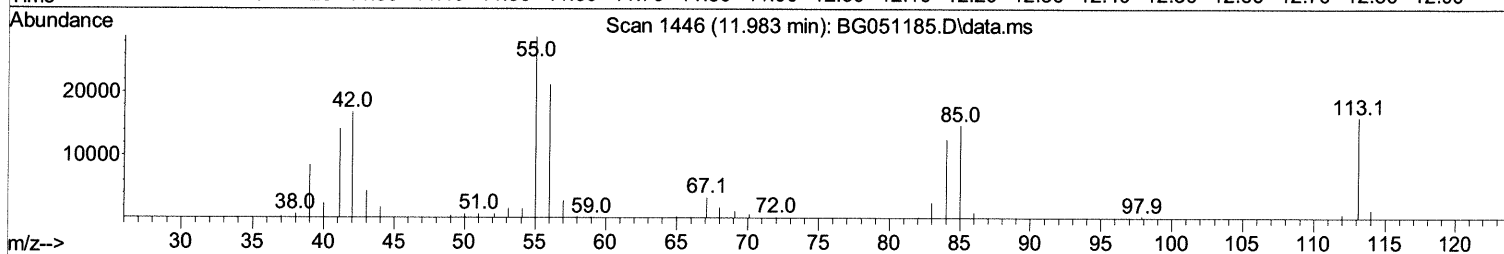
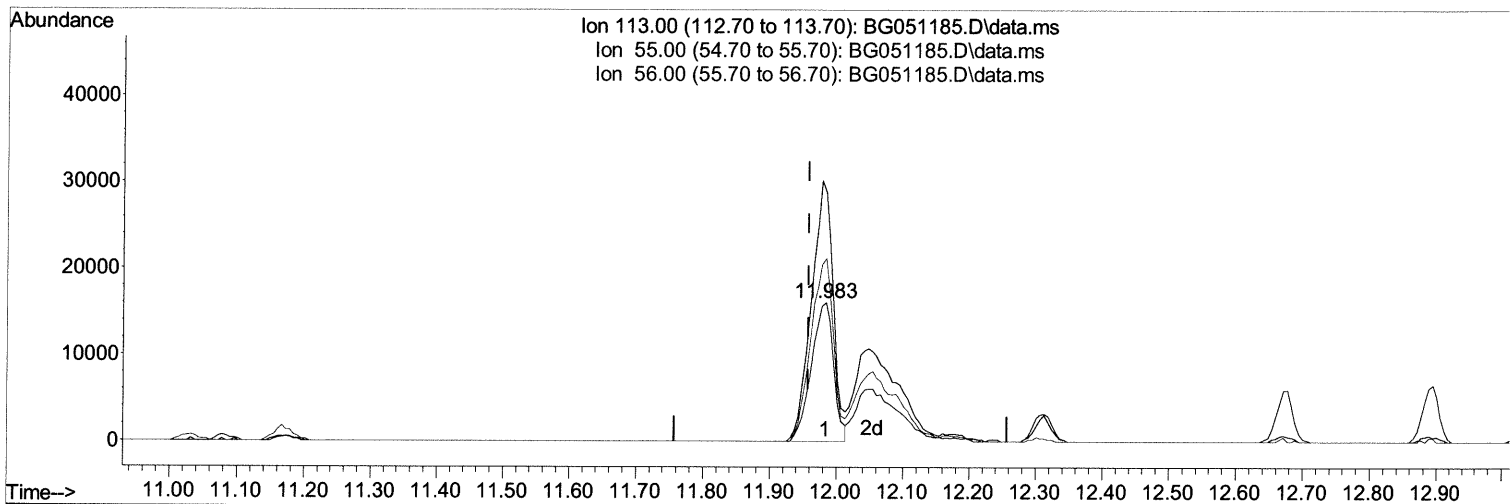
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TIC: BG051185.D\data.ms

(34) Caprolactam

11.983min (+ 0.026) 51.79 ng/ul

response 37043

| Ion | Exp% | Act% |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 183.80 | 179.26 |
| 56.00 | 136.50 | 131.85 |
| 0.00 | 0.00 | 0.00 |

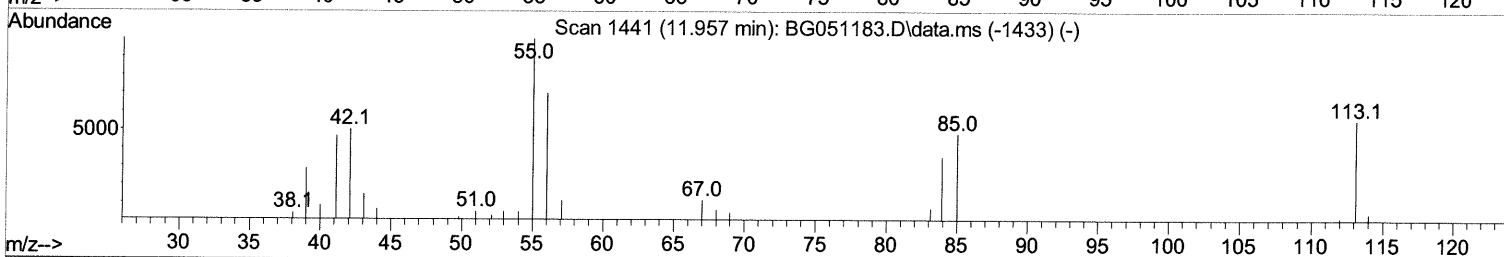
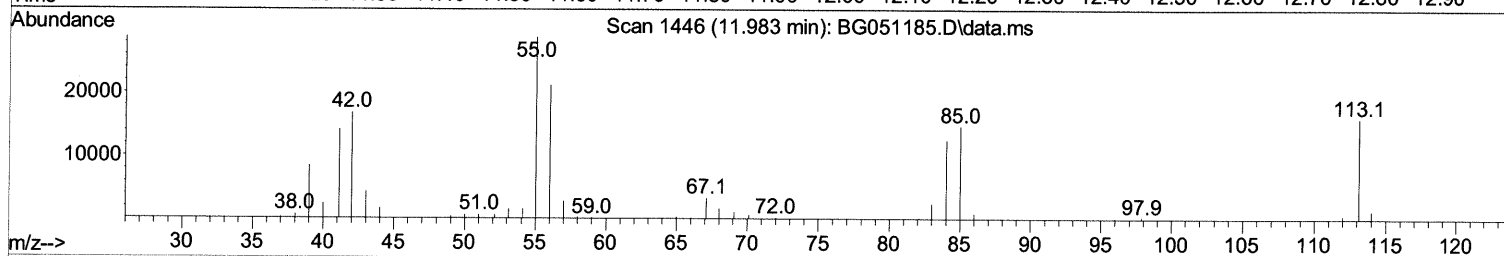
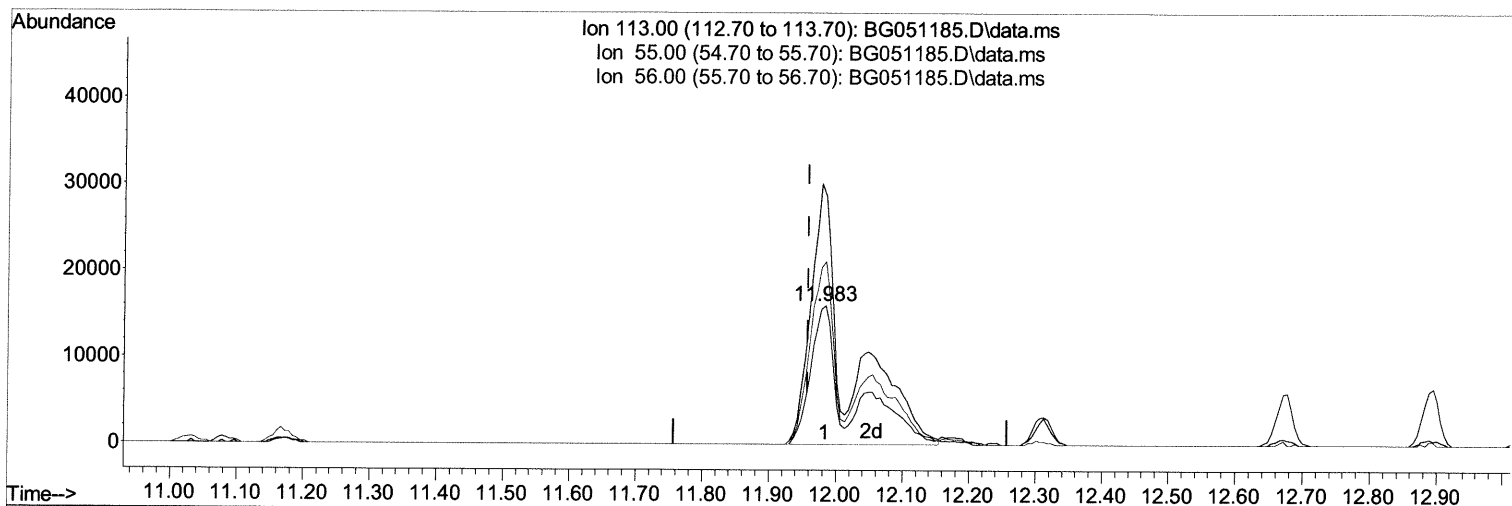
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051185.D
Acq On : 23 Nov 2021 13:37
Operator : CG/JU
Sample : SST08023
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTD080424

Manual IntegrationsAPPROVED

Quant Time: Nov 23 14:56:53 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 23 13:54:17 2021
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/23/2021
Supervised By :mohammad ahmed 11/30/2021



TIC: BG051185.D\data.ms

(34) Caprolactam

11.983min (+ 0.026) 89.26 ng/ul m 11/24/21 JU

response 63844

| Ion | Exp% | Act% |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 183.80 | 179.26 |
| 56.00 | 136.50 | 131.85 |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
 Data File : BG051185.D
 Acq On : 23 Nov 2021 13:37
 Operator : CG/JU
 Sample : SSTD08023
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD080424

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/23/2021
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Quant Time: Nov 23 14:56:53 2021
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 23 13:54:17 2021
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|------------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 8.200 | 152 | 23338 | 20.000 | ng/ul | 0.00 |
| 20) Naphthalene-d8 | 11.032 | 136 | 108519 | 20.000 | ng/ul | 0.00 |
| 38) Acenaphthene-d10 | 14.833 | 164 | 72994 | 20.000 | ng/ul | 0.00 |
| 64) Phenanthrene-d10 | 17.583 | 188 | 164888 | 20.000 | ng/ul | 0.00 |
| 79) Chrysene-d12 | 21.889 | 240 | 138377 | 20.000 | ng/ul | 0.00 |
| 88) Perylene-d12 | 25.291 | 264 | 142201 | 20.000 | ng/ul | 0.00 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.540 | 96 | 25263 | 34.938 | ng/uL | 0.00 |
| 4) Pyridine-d5 | 3.969 | 84 | 185424 | 85.718 | ng/ul | 0.00 |
| 7) Phenol-d5 | 7.359 | 99 | 220744 | 88.661 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.518 | 67 | 135082 | 83.990 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4 | 7.736 | 132 | 160103 | 92.788 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8 | 8.922 | 113 | 176167 | 89.878 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5 | 9.381 | 128 | 85590 | 92.810 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4 | 10.109 | 143 | 100372 | 97.885 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.656 | 165 | 170391 | 98.646 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4 | 11.173 | 131 | 240297 | 91.864 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6 | 14.234 | 166 | 513976 | 92.036 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8 | 14.533 | 160 | 646323 | 92.894 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4 | 15.056 | 143 | 88997 | 87.893 | ng/ul | 0.00 |
| 60) Fluorene-d10 | 15.826 | 176 | 456600 | 92.297 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.961 | 200 | 100972 | 101.005 | ng/ul | 0.00 |
| 73) Anthracene-d10 | 17.689 | 188 | 688210 | 88.281 | ng/ul | 0.00 |
| 81) Pyrene-d10 | 19.962 | 212 | 772058 | 86.383 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12 | 25.062 | 264 | 714320 | 90.868 | ng/ul | 0.02 |
| Target Compounds | | | | | | |
| | | | | | Qvalue | |
| 2) 1,4-Dioxane | 3.576 | 88 | 25775 | 32.454 | ng/uL | 98 |
| 5) Pyridine | 3.987 | 79 | 192847 | 86.123 | ng/ul | 98 |
| 6) Benzaldehyde | 7.336 | 77 | 104652 | 66.642 | ng/ul | 91 |
| 8) Phenol | 7.389 | 94 | 225686 | 87.627 | ng/ul | 99 |
| 10) Bis(2-Chloroethyl)ether | 7.612 | 93 | 168507 | 87.408 | ng/ul | 98 |
| 12) 2-Chlorophenol | 7.765 | 128 | 160640 | 91.694 | ng/ul | 96 |
| 13) 2-Methylphenol | 8.646 | 108 | 171419 | 90.046 | ng/ul | 98 |
| 14) 2,2'-oxybis(1-Chloropr... | 8.723 | 45 | 241157 | 79.419 | ng/ul | 98 |
| 16) Acetophenone | 9.034 | 105 | 266898 | 87.655 | ng/ul | 98 |
| 17) N-Nitroso-di-n-propyla... | 9.016 | 70 | 156516 | 85.196 | ng/ul | 97 |
| 18) 4-Methylphenol | 8.987 | 108 | 180278 | 88.941 | ng/ul | 96 |
| 19) Hexachloroethane | 9.287 | 117 | 66242 | 90.435 | ng/ul | 96 |
| 22) Nitrobenzene | 9.428 | 77 | 223690 | 86.971 | ng/ul | 96 |
| 23) Isophorone | 9.951 | 82 | 436176 | 87.380 | ng/ul | 98 |
| 25) 2-Nitrophenol | 10.139 | 139 | 99948 | 97.156 | ng/ul | 98 |
| 26) 2,4-Dimethylphenol | 10.191 | 107 | 203205 | 89.752 | ng/ul | 100 |
| 27) Bis(2-Chloroethoxy)met... | 10.421 | 93 | 240165 | 89.295 | ng/ul | 99 |
| 29) 2,4-Dichlorophenol | 10.679 | 162 | 161088 | 95.609 | ng/ul | 97 |
| 30) Naphthalene | 11.085 | 128 | 532424 | 89.723 | ng/ul | 99 |
| 32) 4-Chloroaniline | 11.196 | 127 | 240157 | 92.461 | ng/ul | 100 |
| 33) Hexachlorobutadiene | 11.349 | 225 | 107417 | 97.136 | ng/ul | 99 |
| 34) Caprolactam | 11.983 | 113 | 63844m | 89.256 | ng/ul | > 11/29/21 |
| 35) 4-Chloro-3-methylphenol | 12.312 | 107 | 198705 | 92.321 | ng/ul | 98 |

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 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual IntegrationsAPPROVED

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 Quant Title : SVOA CALIBRATION
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| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| 36) 2-Methylnaphthalene | 12.671 | 142 | 366858 | 90.747 | ng/ul | 100 |
| 37) 1-Methylnaphthalene | 12.894 | 142 | 374429 | 91.407 | ng/ul | 98 |
| 39) 1,2,4,5-Tetrachloroben... | 13.035 | 216 | 214150 | 100.688 | ng/ul | 97 |
| 40) Hexachlorocyclopentadiene | 13.006 | 237 | 86836 | 85.021 | ng/ul | 99 |
| 41) 2,4,6-Trichlorophenol | 13.276 | 196 | 142622 | 102.479 | ng/ul | 98 |
| 42) 2,4,5-Trichlorophenol | 13.358 | 196 | 148380 | 99.307 | ng/ul | 98 |
| 43) 1,1'-Biphenyl | 13.670 | 154 | 490289 | 91.894 | ng/ul | 99 |
| 44) 2-Chloronaphthalene | 13.723 | 162 | 390501 | 93.399 | ng/ul | 98 |
| 45) 2-Nitroaniline | 13.928 | 65 | 147304 | 88.677 | ng/ul | 90 |
| 47) Dimethylphthalate | 14.281 | 163 | 507671 | 90.917 | ng/ul | 100 |
| 48) 2,6-Dinitrotoluene | 14.416 | 165 | 115490 | 98.819 | ng/ul | 91 |
| 50) Acenaphthylene | 14.563 | 152 | 619326 | 88.816 | ng/ul | 99 |
| 51) 3-Nitroaniline | 14.751 | 138 | 102810 | 85.056 | ng/ul | 95 |
| 52) Acenaphthene | 14.898 | 153 | 420979 | 91.813 | ng/ul | 96 |
| 53) 2,4-Dinitrophenol | 14.968 | 184 | 64411 | 99.905 | ng/ul | 88 |
| 55) 4-Nitrophenol | 15.074 | 109 | 77875 | 83.860 | ng/ul | 93 |
| 56) Dibenzofuran | 15.233 | 168 | 587184 | 89.467 | ng/ul | 97 |
| 57) 2,4-Dinitrotoluene | 15.209 | 165 | 162036 | 97.150 | ng/ul | 91 |
| 58) 2,3,4,6-Tetrachlorophenol | 15.462 | 232 | 123011 | 104.761 | ng/ul | 96 |
| 59) Diethylphthalate | 15.632 | 149 | 535227 | 89.547 | ng/ul | 97 |
| 61) Fluorene | 15.885 | 166 | 472137 | 90.888 | ng/ul | 100 |
| 62) 4-Chlorophenyl-phenyle... | 15.867 | 204 | 256295 | 94.748 | ng/ul | 95 |
| 63) 4-Nitroaniline | 15.920 | 138 | 98117 | 81.823 | ng/ul | 96 |
| 66) 4,6-Dinitro-2-methylph... | 15.973 | 198 | 96678 | 99.168 | ng/ul | 95 |
| 67) N-Nitrosodiphenylamine | 16.085 | 169 | 427250 | 92.699 | ng/ul | 99 |
| 68) 4-Bromophenyl-phenylether | 16.760 | 248 | 165781 | 101.075 | ng/ul | 93 |
| 69) Hexachlorobenzene | 16.889 | 284 | 167706 | 99.459 | ng/ul | 96 |
| 70) Atrazine | 17.030 | 200 | 181173 | 92.695 | ng/ul | 99 |
| 71) Pentachlorophenol | 17.236 | 266 | 81388 | 105.133 | ng/ul | 96 |
| 72) Phenanthrene | 17.630 | 178 | 801968 | 91.116 | ng/ul | 98 |
| 74) Anthracene | 17.724 | 178 | 779196 | 88.232 | ng/ul | 99 |
| 75) 1,2,3,4-Tetrachloroben... | 13.640 | 216 | 228354 | 101.712 | ng/uL | 97 |
| 76) Pentachlorobenzene | 15.156 | 250 | 209576 | 100.746 | ng/uL | 99 |
| 77) Carbazole | 17.994 | 167 | 710039 | 89.703 | ng/ul | 98 |
| 78) Di-n-butylphthalate | 18.517 | 149 | 892715 | 85.830 | ng/ul | 99 |
| 80) Fluoranthene | 19.633 | 202 | 946643 | 88.256 | ng/ul | 97 |
| 82) Pyrene | 19.998 | 202 | 895658 | 85.459 | ng/ul | 97 |
| 83) Butylbenzylphthalate | 20.855 | 149 | 397713 | 88.247 | ng/ul | 93 |
| 84) 3,3'-Dichlorobenzidine | 21.772 | 252 | 293200 | 87.004 | ng/ul | 99 |
| 85) Benzo(a)anthracene | 21.872 | 228 | 856666 | 89.425 | ng/ul | 97 |
| 86) Bis(2-ethylhexyl)phtha... | 21.725 | 149 | 562931 | 87.017 | ng/ul | 98 |
| 87) Chrysene | 21.936 | 228 | 822637 | 89.893 | ng/ul | 97 |
| 89) Di-n-octyl phthalate | 22.994 | 149 | 953941 | 82.401 | ng/ul | 100 |
| 90) Benzo(b)fluoranthene | 24.210 | 252 | 886111 | 87.471 | ng/ul | 98 |
| 91) Benzo(k)fluoranthene | 24.281 | 252 | 825309 | 86.820 | ng/ul | 98 |
| 93) Benzo(a)pyrene | 25.145 | 252 | 845424 | 87.624 | ng/ul | 97 |
| 94) Indeno(1,2,3-cd)pyrene | 29.222 | 276 | 972330 | 90.530 | ng/ul | 98 |
| 95) Dibenzo(a,h)anthracene | 29.287 | 278 | 810855 | 89.225 | ng/ul | 98 |
| 96) Benzo(g,h,i)perylene | 30.468 | 276 | 810492 | 90.149 | ng/ul | 96 |

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