

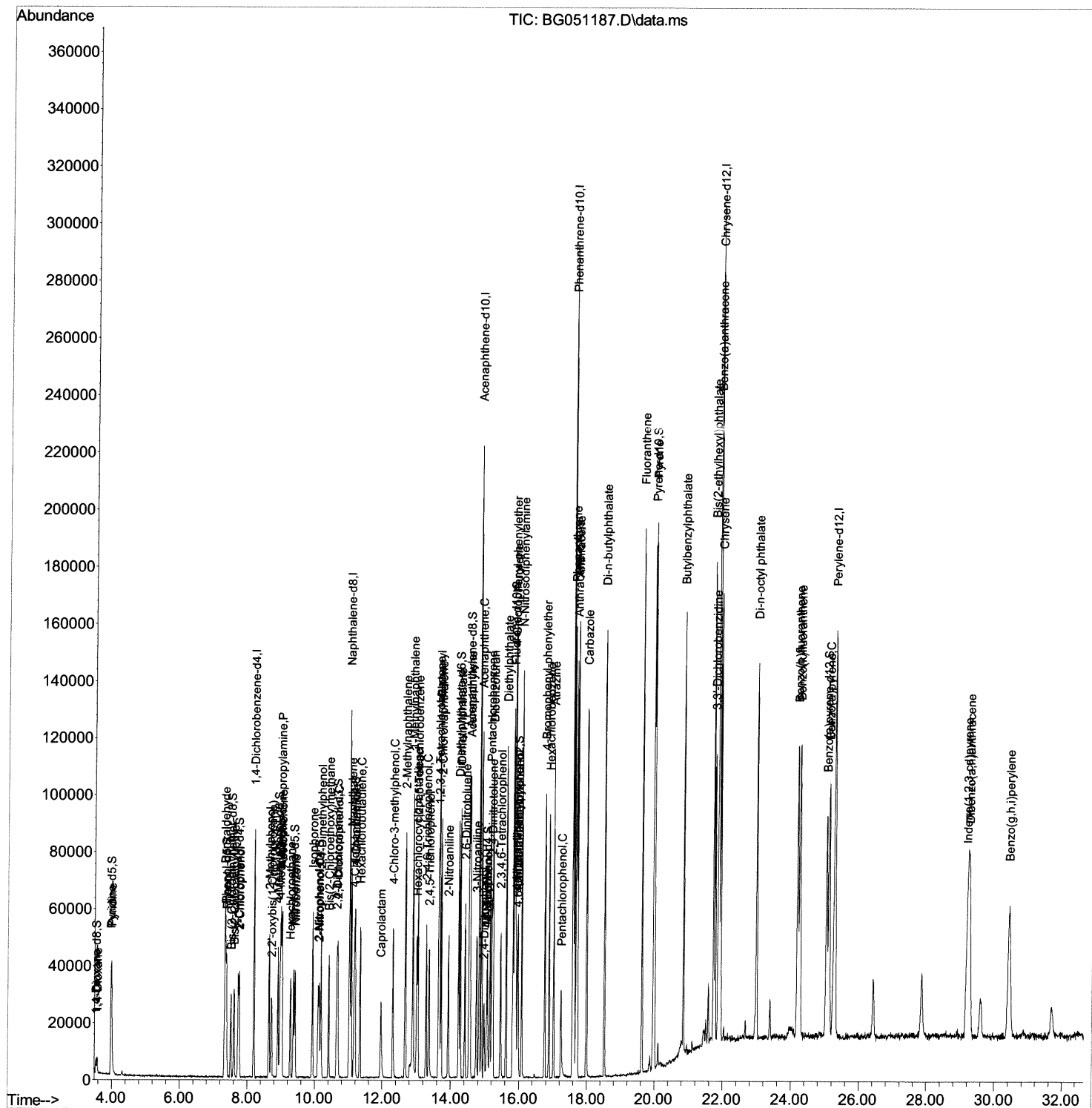
```
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\  
Data File : BG051187.D  
Acq On    : 23 Nov 2021  15:58  
Operator  : CG/JU  
Sample    : SST001020  
Misc      :  
ALS Vial  : 8    Sample Multiplier: 1
```

Instrument :
BNA_G
ClientSampleId :
SSTD010426

Manual IntegrationsAPPROVED

Quant Time: Nov 23 16:42:16 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 23 15:29:22 2021
Response via : Initial Calibration

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



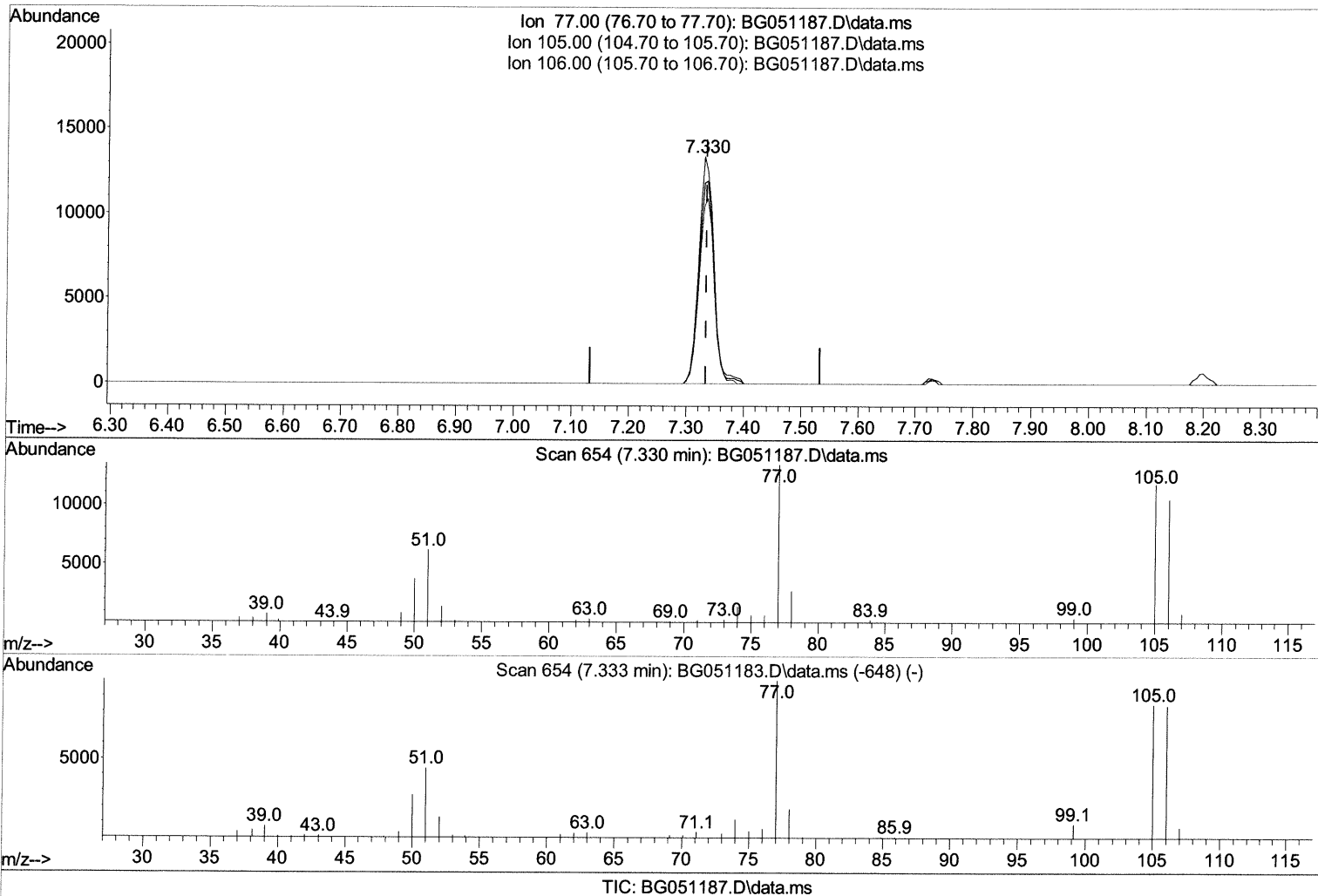
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051187.D
Acq On : 23 Nov 2021 15:58
Operator : CG/JU
Sample : SSTD01020
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTD010426

Manual IntegrationsAPPROVED

Quant Time: Nov 23 16:42:16 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 23 15:29:22 2021
Response via : Initial Calibration

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



(6) Benzaldehyde

7.330min (-0.004) 17.31 ng/ul

response 24448

| Ion | Exp% | Act% |
|--------|--------|--------|
| 77.00 | 100.00 | 100.00 |
| 105.00 | 88.00 | 88.18 |
| 106.00 | 76.50 | 78.44 |
| 0.00 | 0.00 | 0.00 |

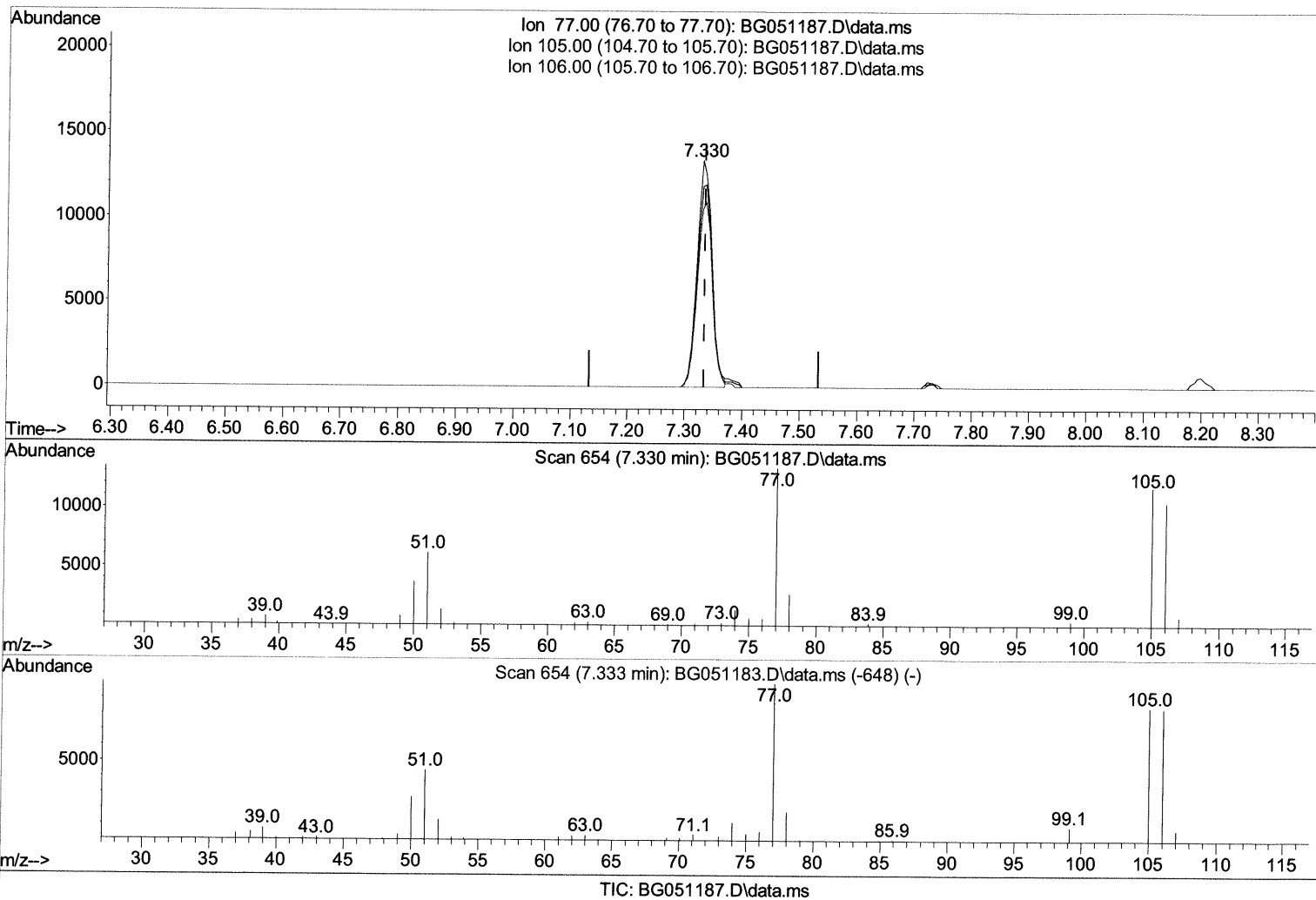
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
Data File : BG051187.D
Acq On : 23 Nov 2021 15:58
Operator : CG/JU
Sample : SSTD01020
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTD010426

Manual IntegrationsAPPROVED

Quant Time: Nov 23 16:42:16 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 23 15:29:22 2021
Response via : Initial Calibration

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



(6) Benzaldehyde

7.330min (-0.004) 16.94 ng/ul m 11/29/21JU

response 23932

| Ion | Exp% | Act% |
|--------|--------|--------|
| 77.00 | 100.00 | 100.00 |
| 105.00 | 88.00 | 88.18 |
| 106.00 | 76.50 | 78.44 |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
 Data File : BG051187.D
 Acq On : 23 Nov 2021 15:58
 Operator : CG/JU
 Sample : SST01020
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SST010426

Manual Integrations APPROVED

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

Quant Time: Nov 23 16:42:16 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 23 15:29:22 2021
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 8.199 | 152 | 24591 | 20.000 | ng/ul | 0.00 |
| 20) Naphthalene-d8 | 11.025 | 136 | 110381 | 20.000 | ng/ul | 0.00 |
| 38) Acenaphthene-d10 | 14.833 | 164 | 78299 | 20.000 | ng/ul | 0.00 |
| 64) Phenanthrene-d10 | 17.582 | 188 | 178109 | 20.000 | ng/ul | 0.00 |
| 79) Chrysene-d12 | 21.889 | 240 | 162571 | 20.000 | ng/ul | 0.00 |
| 88) Perylene-d12 | 25.285 | 264 | 159934 | 20.000 | ng/ul | 0.00 |

| | | | | | | |
|-------------------------------|--------|-----|-------|--------|-------|------|
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.534 | 96 | 2772 | 3.817 | ng/uL | 0.00 |
| 4) Pyridine-d5 | 3.969 | 84 | 20004 | 9.647 | ng/ul | 0.00 |
| 7) Phenol-d5 | 7.359 | 99 | 22530 | 9.264 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.512 | 67 | 15329 | 10.169 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4 | 7.729 | 132 | 17307 | 9.952 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8 | 8.910 | 113 | 18706 | 9.699 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5 | 9.374 | 128 | 9653 | 10.512 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4 | 10.103 | 143 | 10330 | 9.796 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.655 | 165 | 17494 | 9.851 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4 | 11.166 | 131 | 25930 | 10.019 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6 | 14.227 | 166 | 61347 | 10.264 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8 | 14.533 | 160 | 76971 | 10.150 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4 | 15.050 | 143 | 8506 | 8.859 | ng/ul | 0.00 |
| 60) Fluorene-d10 | 15.826 | 176 | 55497 | 10.316 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.949 | 200 | 9252 | 8.436 | ng/ul | 0.00 |
| 73) Anthracene-d10 | 17.682 | 188 | 88567 | 10.451 | ng/ul | 0.00 |
| 81) Pyrene-d10 | 19.962 | 212 | 97784 | 9.925 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12 | 25.050 | 264 | 84975 | 9.964 | ng/ul | 0.00 |

| | | | | | | |
|-------------------------------|--------|-----|--------|--------------|----|-------------|
| Target Compounds | | | | Qvalue | | |
| 2) 1,4-Dioxane | 3.575 | 88 | 3151 | 3.825 ng/uL# | 95 | |
| 5) Pyridine | 3.992 | 79 | 20684 | 9.505 ng/ul | 95 | |
| 6) Benzaldehyde | 7.330 | 77 | 23932m | 16.944 ng/ul | > | 11/29/21 JU |
| 8) Phenol | 7.382 | 94 | 24372 | 9.692 ng/ul | 97 | |
| 10) Bis(2-Chloroethyl)ether | 7.606 | 93 | 18858 | 9.895 ng/ul | 94 | |
| 12) 2-Chlorophenol | 7.764 | 128 | 17320 | 9.676 ng/ul | 94 | |
| 13) 2-Methylphenol | 8.646 | 108 | 17288 | 9.262 ng/ul | 96 | |
| 14) 2,2'-oxybis(1-Chloropr... | 8.722 | 45 | 28084 | 10.312 ng/ul | 97 | |
| 16) Acetophenone | 9.028 | 105 | 31307 | 10.459 ng/ul | 97 | |
| 17) N-Nitroso-di-n-propyla... | 8.998 | 70 | 17936 | 10.466 ng/ul | 96 | |
| 18) 4-Methylphenol | 8.975 | 108 | 19546 | 9.825 ng/ul | 93 | |
| 19) Hexachloroethane | 9.280 | 117 | 7538 | 10.024 ng/ul | 94 | |
| 22) Nitrobenzene | 9.421 | 77 | 24583 | 10.015 ng/ul | 95 | |
| 23) Isophorone | 9.932 | 82 | 48352 | 10.241 ng/ul | 98 | |
| 25) 2-Nitrophenol | 10.138 | 139 | 11110 | 10.231 ng/ul | 99 | |
| 26) 2,4-Dimethylphenol | 10.185 | 107 | 21645 | 9.663 ng/ul | 94 | |
| 27) Bis(2-Chloroethoxy)met... | 10.414 | 93 | 26873 | 10.279 ng/ul | 96 | |
| 29) 2,4-Dichlorophenol | 10.679 | 162 | 17267 | 9.836 ng/ul | 94 | |
| 30) Naphthalene | 11.078 | 128 | 60691 | 10.085 ng/ul | 98 | |
| 32) 4-Chloroaniline | 11.190 | 127 | 26296 | 10.083 ng/ul | 99 | |
| 33) Hexachlorobutadiene | 11.343 | 225 | 12319 | 10.217 ng/ul | 96 | |
| 34) Caprolactam | 11.954 | 113 | 6899 | 10.222 ng/ul | 95 | |
| 35) 4-Chloro-3-methylphenol | 12.306 | 107 | 21443 | 10.193 ng/ul | 94 | |

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
 Data File : BG051187.D
 Acq On : 23 Nov 2021 15:58
 Operator : CG/JU
 Sample : SSTD01020
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD010426

Manual IntegrationsAPPROVED

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

Quant Time: Nov 23 16:42:16 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 23 15:29:22 2021
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene | 12.670 | 142 | 42821 | 10.611 | ng/ul | 99 |
| 37) 1-Methylnaphthalene | 12.888 | 142 | 43034 | 10.362 | ng/ul | 97 |
| 39) 1,2,4,5-Tetrachloroben... | 13.035 | 216 | 24275 | 9.848 | ng/ul | 96 |
| 40) Hexachlorocyclopentadiene | 12.999 | 237 | 14145 | 18.410 | ng/ul# | 98 |
| 41) 2,4,6-Trichlorophenol | 13.281 | 196 | 14504 | 9.393 | ng/ul | 97 |
| 42) 2,4,5-Trichlorophenol | 13.364 | 196 | 15992 | 9.915 | ng/ul | 97 |
| 43) 1,1'-Biphenyl | 13.669 | 154 | 59293 | 10.118 | ng/ul | 98 |
| 44) 2-Chloronaphthalene | 13.716 | 162 | 47530 | 10.238 | ng/ul | 99 |
| 45) 2-Nitroaniline | 13.922 | 65 | 16107 | 10.118 | ng/ul | 92 |
| 47) Dimethylphthalate | 14.274 | 163 | 62824 | 10.425 | ng/ul | 99 |
| 48) 2,6-Dinitrotoluene | 14.410 | 165 | 12705 | 10.059 | ng/ul | 92 |
| 50) Acenaphthylene | 14.562 | 152 | 77706 | 10.426 | ng/ul | 97 |
| 51) 3-Nitroaniline | 14.744 | 138 | 13982 | 11.311 | ng/ul | 99 |
| 52) Acenaphthene | 14.897 | 153 | 50544 | 10.265 | ng/ul | 99 |
| 53) 2,4-Dinitrophenol | 14.968 | 184 | 8391 | 14.126 | ng/ul# | 86 |
| 55) 4-Nitrophenol | 15.068 | 109 | 7161 | 8.543 | ng/ul | 85 |
| 56) Dibenzofuran | 15.232 | 168 | 72994 | 10.284 | ng/ul | 98 |
| 57) 2,4-Dinitrotoluene | 15.203 | 165 | 18521 | 10.228 | ng/ul | 93 |
| 58) 2,3,4,6-Tetrachlorophenol | 15.461 | 232 | 12192 | 9.732 | ng/ul | 97 |
| 59) Diethylphthalate | 15.626 | 149 | 64941 | 10.210 | ng/ul | 98 |
| 61) Fluorene | 15.878 | 166 | 59085 | 10.380 | ng/ul | 95 |
| 62) 4-Chlorophenyl-phenyle... | 15.861 | 204 | 31891 | 10.465 | ng/ul | 91 |
| 63) 4-Nitroaniline | 15.908 | 138 | 13205 | 10.663 | ng/ul | 97 |
| 66) 4,6-Dinitro-2-methylph... | 15.967 | 198 | 9140 | 8.660 | ng/ul# | 87 |
| 67) N-Nitrosodiphenylamine | 16.078 | 169 | 52690 | 10.355 | ng/ul | 99 |
| 68) 4-Bromophenyl-phenylether | 16.760 | 248 | 19241 | 10.096 | ng/ul | 95 |
| 69) Hexachlorobenzene | 16.883 | 284 | 19343 | 9.969 | ng/ul | 93 |
| 70) Atrazine | 17.018 | 200 | 21828 | 10.205 | ng/ul | 95 |
| 71) Pentachlorophenol | 17.241 | 266 | 9403 | 12.421 | ng/ul | 93 |
| 72) Phenanthrene | 17.623 | 178 | 101712 | 10.404 | ng/ul | 98 |
| 74) Anthracene | 17.717 | 178 | 101680 | 10.402 | ng/ul | 98 |
| 75) 1,2,3,4-Tetrachloroben... | 13.640 | 216 | 25714 | 9.874 | ng/ul | 98 |
| 76) Pentachlorobenzene | 15.150 | 250 | 24593 | 10.191 | ng/ul | 95 |
| 77) Carbazole | 17.988 | 167 | 90021 | 10.530 | ng/ul | 98 |
| 78) Di-n-butylphthalate | 18.516 | 149 | 112838 | 10.192 | ng/ul | 99 |
| 80) Fluoranthene | 19.627 | 202 | 121606 | 10.098 | ng/ul | 97 |
| 82) Pyrene | 19.991 | 202 | 121973 | 10.328 | ng/ul | 97 |
| 83) Butylbenzylphthalate | 20.855 | 149 | 48380 | 9.859 | ng/ul | 94 |
| 84) 3,3'-Dichlorobenzidine | 21.771 | 252 | 38572 | 10.065 | ng/ul | 97 |
| 85) Benzo(a)anthracene | 21.866 | 228 | 111203 | 10.069 | ng/ul | 99 |
| 86) Bis(2-ethylhexyl)phtha... | 21.724 | 149 | 70869 | 10.069 | ng/ul | 98 |
| 87) Chrysene | 21.936 | 228 | 106392 | 10.024 | ng/ul | 99 |
| 89) Di-n-octyl phthalate | 22.994 | 149 | 118314 | 10.243 | ng/ul | 100 |
| 90) Benzo(b)fluoranthene | 24.192 | 252 | 109032 | 10.093 | ng/ul | 96 |
| 91) Benzo(k)fluoranthene | 24.269 | 252 | 102094 | 10.102 | ng/ul | 98 |
| 93) Benzo(a)pyrene | 25.121 | 252 | 103679 | 10.072 | ng/ul | 99 |
| 94) Indeno(1,2,3-cd)pyrene | 29.192 | 276 | 112895 | 9.776 | ng/ul | 96 |
| 95) Dibenzo(a,h)anthracene | 29.257 | 278 | 96981 | 9.925 | ng/ul# | 96 |
| 96) Benzo(g,h,i)perylene | 30.426 | 276 | 95922 | 9.913 | ng/ul | 97 |

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