

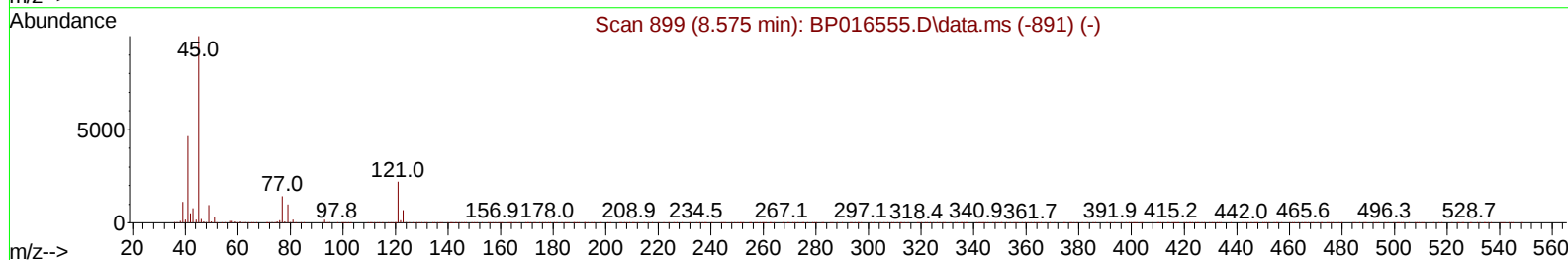
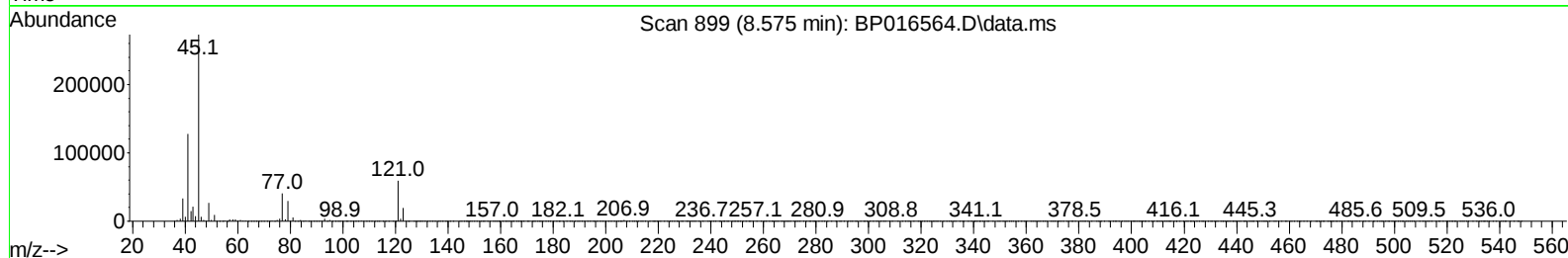
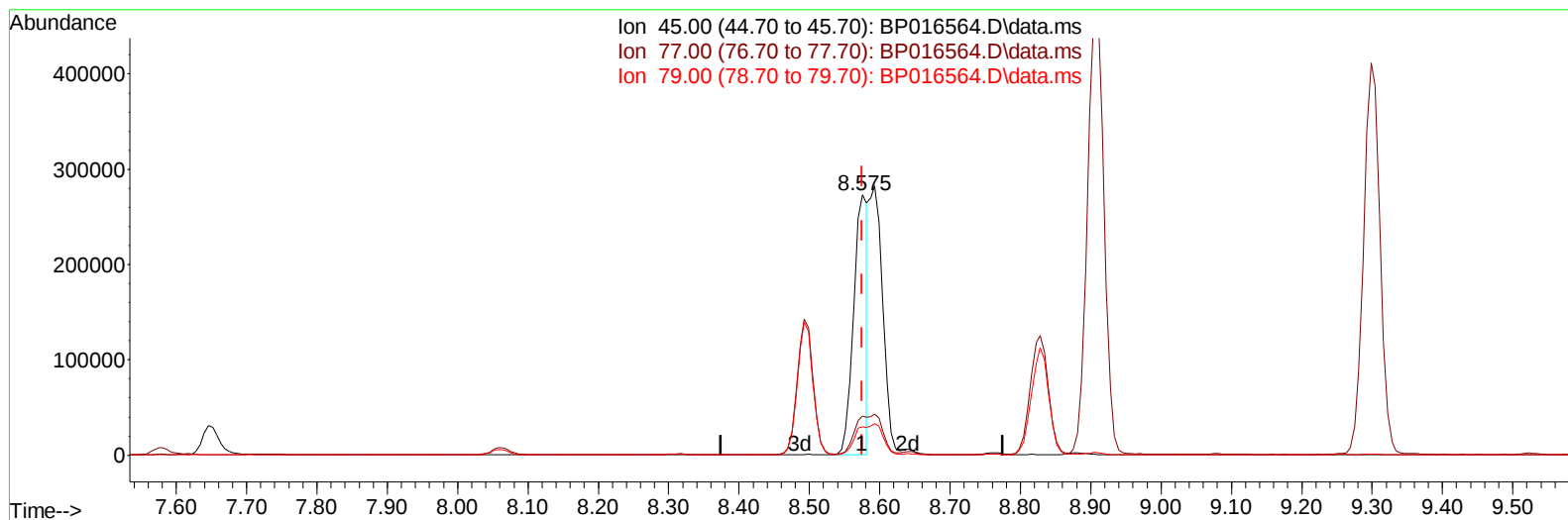
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081423\  
 Data File : BP016564.D  
 Acq On : 14 Aug 2023 16:42  
 Operator : MA/JU  
 Sample : 03965-02MS  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A4832MS

Manual IntegrationsAPPROVED

Quant Time: Aug 14 18:41:18 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 14 18:26:05 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/15/2023  
 Supervised By :Sohil Jodhani 08/15/2023



TIC: BP016564.D\data.ms

(14) 2,2'-oxybis(1-Chloropropane)

8.575min (-0.000) 10.76 ng/u1

response 373377

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 15.10  | 14.92  |
| 79.00 | 10.30  | 10.96  |
| 0.00  | 0.00   | 0.00   |

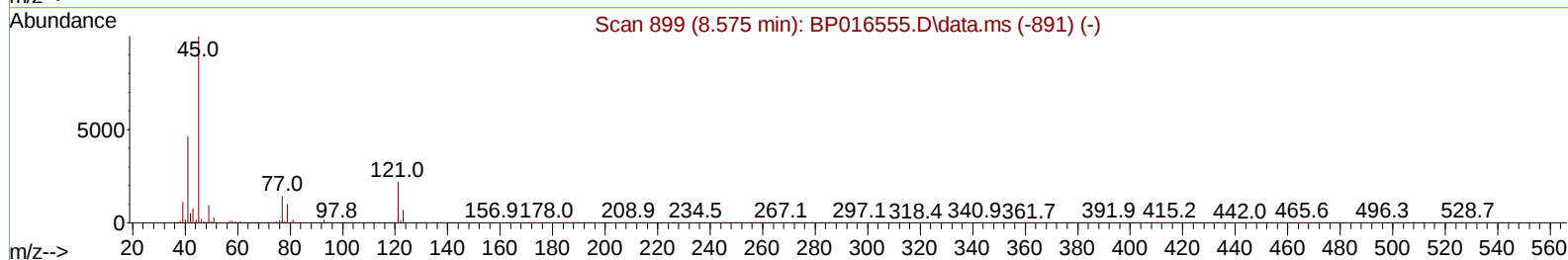
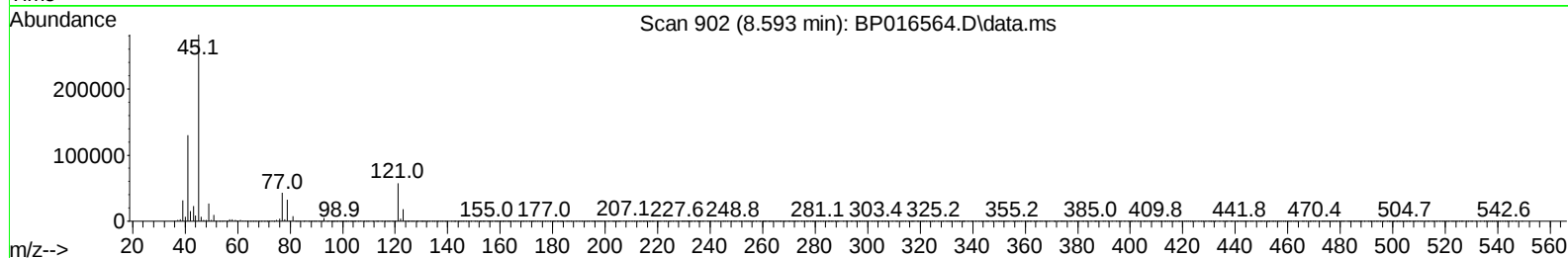
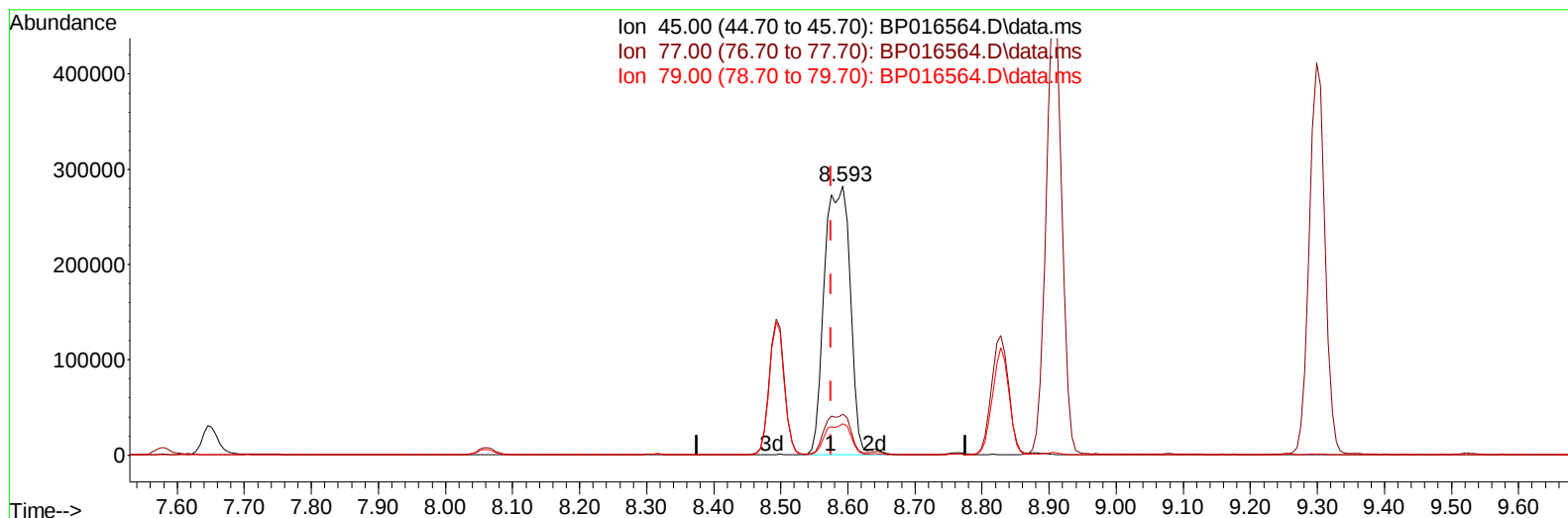
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081423\  
 Data File : BP016564.D  
 Acq On : 14 Aug 2023 16:42  
 Operator : MA/JU  
 Sample : 03965-02MS  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
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 ClientSampleId :  
 A4832MS

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/15/2023  
 Supervised By :Sohil Jodhani 08/15/2023

Quant Time: Aug 14 18:41:51 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 14 18:26:05 2023  
 Response via : Initial Calibration



TIC: BP016564.D\data.ms

## (14) 2,2'-oxybis(1-Chloropropane)

8.593min (+ 0.018) 21.85 ng/ul m

response 758010

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 15.10  | 15.18  |
| 79.00 | 10.30  | 11.60  |
| 0.00  | 0.00   | 0.00   |

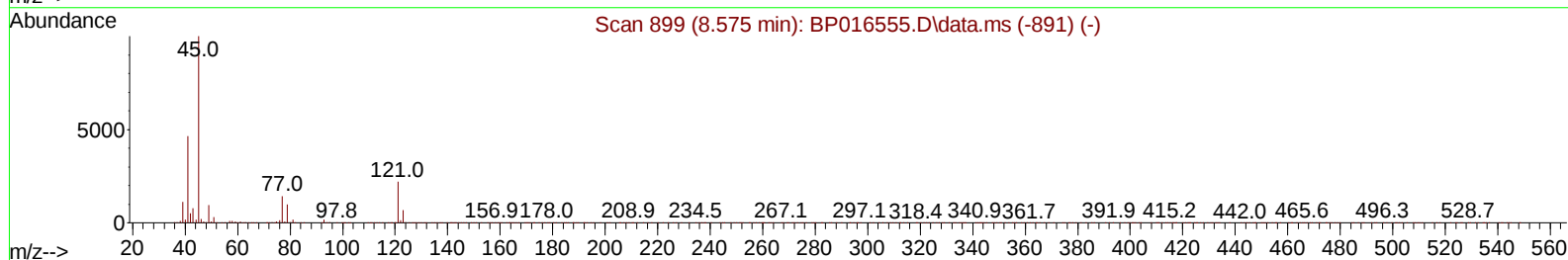
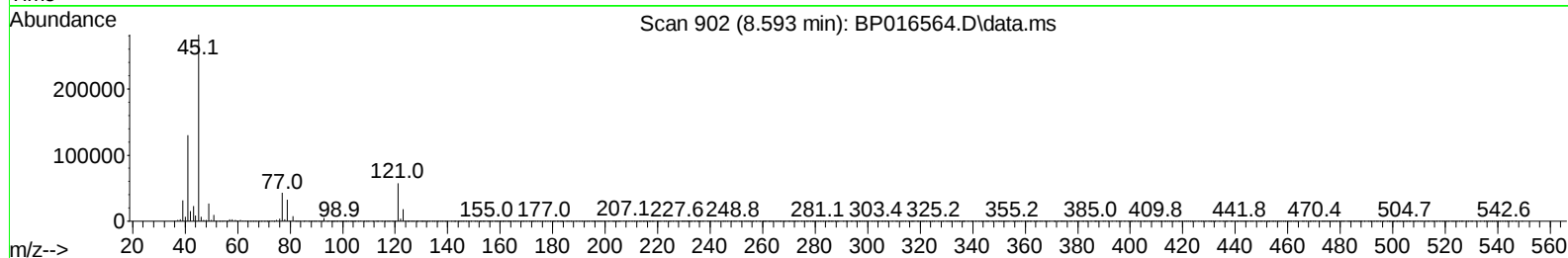
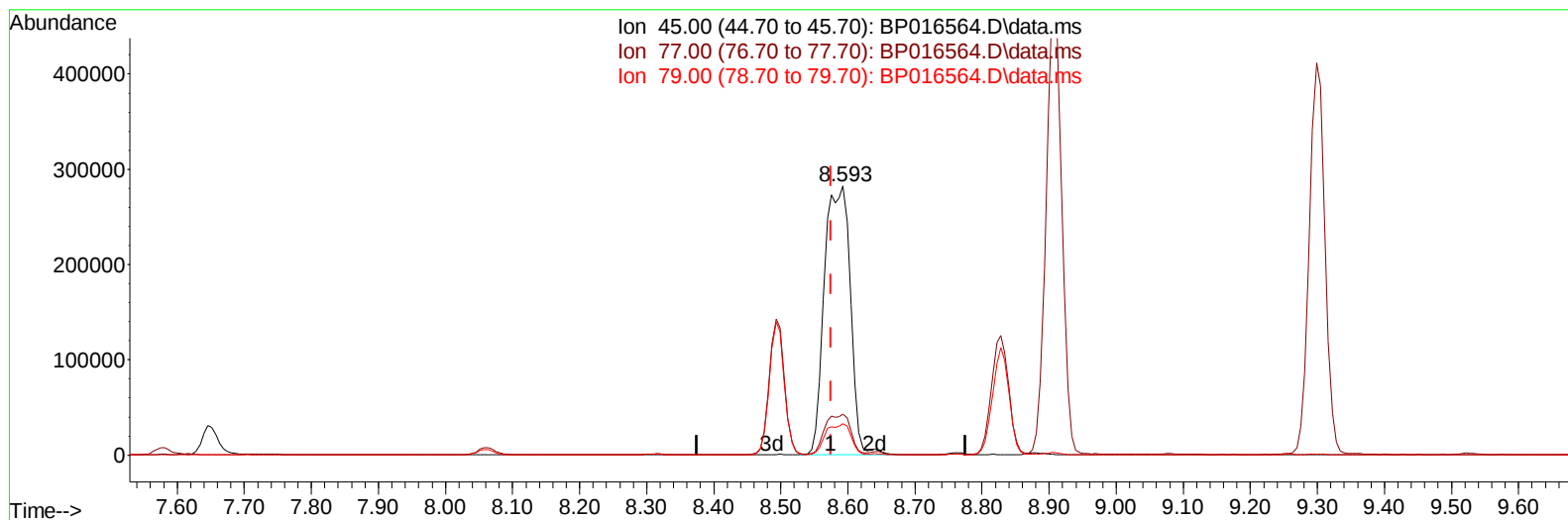
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081423\  
Data File : BP016564.D  
Acq On : 14 Aug 2023 16:42  
Operator : MA/JU  
Sample : 03965-02MS  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A4832MS

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/15/2023  
Supervised By :Sohil Jodhani 08/15/2023

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Quant Title : SVOA CALIBRATION  
QLast Update : Mon Aug 14 18:26:05 2023  
Response via : Initial Calibration



TIC: BP016564.D\data.ms

## (14) 2,2'-oxybis(1-Chloropropane)

8.593min (+ 0.018) 21.85 ng/ul m

response 758010

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 15.10  | 15.18  |
| 79.00 | 10.30  | 11.60  |
| 0.00  | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081423\  
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Manual IntegrationsAPPROVED

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 Quant Title : SVOA CALIBRATION  
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Reviewed By :Yogesh Patel 08/15/2023  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.063  | 152  | 312758   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.893 | 136  | 1414467  | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.710 | 164  | 865160   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.475 | 188  | 1933566  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.574 | 240  | 1917307  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.163 | 264  | 2165405  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.393  | 96   | 28872    | 3.508  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.834  | 84   | 279847   | 11.175 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.199  | 99   | 575331   | 19.953 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.393  | 67   | 367584   | 19.355 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.575  | 132  | 413253   | 19.954 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.763  | 113  | 435735   | 19.005 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.257  | 128  | 203904   | 20.487 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.975  | 143  | 197256   | 20.867 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.499 | 165  | 400023   | 21.504 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.046 | 131  | 416304   | 13.043 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.122 | 166  | 1263778  | 20.196 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.410 | 160  | 1448753  | 20.512 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.904 | 143  | 237689   | 18.646 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.704 | 176  | 1097278  | 20.569 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.822 | 200  | 117194   | 12.740 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.575 | 188  | 1731853  | 20.707 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.810 | 212  | 2136222  | 21.264 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.992 | 264  | 2216196  | 21.356 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.434  | 88   | 81606    | 9.017  | ng/uL  | 98       |
| 5) Pyridine                   | 3.852  | 79   | 352459   | 13.566 | ng/ul  | 98       |
| 6) Benzaldehyde               | 7.210  | 77   | 412482   | 25.951 | ng/ul  | 98       |
| 8) Phenol                     | 7.222  | 94   | 751771   | 24.394 | ng/ul  | 98       |
| 10) Bis(2-Chloroethyl)ether   | 7.493  | 93   | 565669   | 22.723 | ng/ul  | 98       |
| 12) 2-Chlorophenol            | 7.610  | 128  | 512267   | 23.508 | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.493  | 108  | 510914   | 22.824 | ng/ul  | 97       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.593  | 45   | 758010m  | 21.848 | ng/ul  |          |
| 16) Acetophenone              | 8.905  | 105  | 959311   | 25.911 | ng/ul  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.881  | 70   | 461437   | 23.239 | ng/ul  | 99       |
| 18) 4-Methylphenol            | 8.828  | 108  | 570998   | 23.209 | ng/ul  | 98       |
| 19) Hexachloroethane          | 9.134  | 117  | 208315   | 22.882 | ng/ul  | 95       |
| 22) Nitrobenzene              | 9.299  | 77   | 673740   | 24.182 | ng/ul  | 99       |
| 23) Isophorone                | 9.816  | 82   | 1275555  | 24.239 | ng/ul  | 100      |
| 25) 2-Nitrophenol             | 10.004 | 139  | 272306   | 25.198 | ng/ul  | 100      |
| 26) 2,4-Dimethylphenol        | 10.040 | 107  | 529423   | 20.781 | ng/ul  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 10.304 | 93   | 784063   | 23.243 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.528 | 162  | 476387   | 24.890 | ng/ul  | 97       |
| 30) Naphthalene               | 10.946 | 128  | 1733822  | 23.461 | ng/ul  | 100      |
| 32) 4-Chloroaniline           | 11.069 | 127  | 512795   | 16.221 | ng/ul  | 99       |
| 33) Hexachlorobutadiene       | 11.181 | 225  | 269799   | 23.384 | ng/ul  | 100      |
| 34) Caprolactam               | 11.881 | 113  | 161350   | 21.669 | ng/ul  | 95       |
| 35) 4-Chloro-3-methylphenol   | 12.163 | 107  | 587651   | 25.118 | ng/ul  | 98       |

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## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/15/2023  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 14 18:26:05 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.540 | 142  | 1149726  | 23.225 | ng/ul | 99       |
| 37) 1-Methylnaphthalene       | 12.763 | 142  | 1152327  | 23.118 | ng/ul | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.887 | 216  | 549496   | 23.746 | ng/ul | 99       |
| 40) Hexachlorocyclopentadiene | 12.845 | 237  | 277313   | 19.696 | ng/ul | 99       |
| 41) 2,4,6-Trichlorophenol     | 13.140 | 196  | 362653   | 25.181 | ng/ul | 99       |
| 42) 2,4,5-Trichlorophenol     | 13.204 | 196  | 406311   | 26.027 | ng/ul | 98       |
| 43) 1,1'-Biphenyl             | 13.545 | 154  | 1530487  | 23.799 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.592 | 162  | 1188363  | 23.751 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.822 | 65   | 398433   | 26.336 | ng/ul | 94       |
| 47) Dimethylphthalate         | 14.169 | 163  | 1559765  | 24.379 | ng/ul | 99       |
| 48) 2,6-Dinitrotoluene        | 14.310 | 165  | 295600   | 26.674 | ng/ul | 100      |
| 50) Acenaphthylene            | 14.440 | 152  | 1904042  | 22.890 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.645 | 138  | 323062   | 25.999 | ng/ul | 98       |
| 52) Acenaphthene              | 14.775 | 153  | 1327868  | 23.848 | ng/ul | 97       |
| 53) 2,4-Dinitrophenol         | 14.839 | 184  | 92382    | 14.974 | ng/ul | 98       |
| 55) 4-Nitrophenol             | 14.922 | 109  | 264484   | 25.521 | ng/ul | 99       |
| 56) Dibenzofuran              | 15.110 | 168  | 1818651  | 23.939 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 15.087 | 165  | 442355   | 26.656 | ng/ul | 99       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.322 | 232  | 334656   | 25.586 | ng/ul | 99       |
| 59) Diethylphthalate          | 15.522 | 149  | 1627319  | 24.876 | ng/ul | 99       |
| 61) Fluorene                  | 15.763 | 166  | 1516774  | 24.349 | ng/ul | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.751 | 204  | 696369   | 23.814 | ng/ul | 99       |
| 63) 4-Nitroaniline            | 15.810 | 138  | 340114   | 28.602 | ng/ul | 97       |
| 66) 4,6-Dinitro-2-methylph... | 15.834 | 198  | 159658   | 15.654 | ng/ul | 96       |
| 67) N-Nitrosodiphenylamine    | 15.975 | 169  | 1274186  | 24.311 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.657 | 248  | 423434   | 24.527 | ng/ul | 97       |
| 69) Hexachlorobenzene         | 16.739 | 284  | 481699   | 23.879 | ng/ul | 99       |
| 70) Atrazine                  | 16.928 | 200  | 380426   | 20.175 | ng/ul | 98       |
| 71) Pentachlorophenol         | 17.098 | 266  | 242418   | 19.519 | ng/ul | 98       |
| 72) Phenanthrene              | 17.522 | 178  | 2529741  | 24.833 | ng/ul | 100      |
| 74) Anthracene                | 17.616 | 178  | 2485939  | 24.295 | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.498 | 216  | 535106   | 21.776 | ng/uL | 96       |
| 76) Pentachlorobenzene        | 15.004 | 250  | 528343   | 23.963 | ng/uL | 98       |
| 77) Carbazole                 | 17.892 | 167  | 2388103  | 25.326 | ng/ul | 99       |
| 78) Di-n-butylphthalate       | 18.404 | 149  | 3012252  | 27.596 | ng/ul | 100      |
| 80) Fluoranthene              | 19.480 | 202  | 3132353  | 26.124 | ng/ul | 99       |
| 82) Pyrene                    | 19.839 | 202  | 3263464  | 25.775 | ng/ul | 97       |
| 83) Butylbenzylphthalate      | 20.692 | 149  | 1426553  | 27.968 | ng/ul | 97       |
| 84) 3,3'-Dichlorobenzidine    | 21.492 | 252  | 598013   | 14.324 | ng/ul | 98       |
| 85) Benzo(a)anthracene        | 21.557 | 228  | 3224379  | 25.023 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.433 | 149  | 2146631  | 28.999 | ng/ul | 100      |
| 87) Chrysene                  | 21.616 | 228  | 3053592  | 25.325 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 22.427 | 149  | 3649323  | 27.009 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.357 | 252  | 3325346  | 25.791 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 23.410 | 252  | 3287834  | 25.704 | ng/ul | 99       |
| 93) Benzo(a)pyrene            | 24.045 | 252  | 2878377  | 23.666 | ng/ul | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 26.915 | 276  | 3674797  | 25.600 | ng/ul | 99       |
| 95) Dibenzo(a,h)anthracene    | 26.945 | 278  | 3036499  | 25.418 | ng/ul | 99       |
| 96) Benzo(g,h,i)perylene      | 27.774 | 276  | 2691182  | 23.556 | ng/ul | 99       |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

A4832MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 08/15/2023  
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