

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM020623\  
 Data File : BM038697.D  
 Acq On : 06 Feb 2023 17:54  
 Operator : CG/JU  
 Sample : SST04036  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST040036

Quant Time: Feb 07 00:53:10 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM020623.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 07 00:48:02 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 02/07/2023  
 Supervised By :mohammad ahmed 02/07/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.904  | 152  | 79408    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.716 | 136  | 325948   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.551 | 164  | 264026   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.292 | 188  | 567684   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.468 | 240  | 472674   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.850 | 264  | 491575   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.281  | 96   | 22345    | 10.033 | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.710  | 84   | 151001   | 27.077 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.087  | 99   | 204960   | 38.668 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.245  | 67   | 113875   | 36.806 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.440  | 132  | 197534   | 44.938 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.639  | 113  | 186429   | 44.081 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.086  | 128  | 101256   | 40.983 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.804  | 143  | 138256   | 42.373 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.345 | 165  | 278359   | 38.548 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.869 | 131  | 300986   | 42.890 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.969 | 166  | 855674   | 39.233 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.245 | 160  | 969107   | 41.755 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.780 | 143  | 140210   | 46.471 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.539 | 176  | 726173   | 37.231 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.674 | 200  | 140053   | 28.821 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.392 | 188  | 1089323  | 38.122 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.686 | 212  | 1124163  | 46.156 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.703 | 264  | 1054510  | 41.793 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.316  | 88   | 22717    | 9.467  | ng/uL  | 99       |
| 5) Pyridine                   | 3.728  | 79   | 156117   | 27.259 | ng/u1  | 98       |
| 6) Benzaldehyde               | 7.051  | 77   | 95942    | 39.256 | ng/u1  | 95       |
| 8) Phenol                     | 7.116  | 94   | 203802   | 38.403 | ng/u1  | 96       |
| 10) Bis(2-Chloroethyl)ether   | 7.340  | 93   | 165298   | 43.594 | ng/u1  | 99       |
| 12) 2-Chlorophenol            | 7.475  | 128  | 191438   | 43.401 | ng/u1  | 97       |
| 13) 2-Methylphenol            | 8.369  | 108  | 171461   | 42.524 | ng/u1  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.445  | 45   | 146783   | 37.469 | ng/u1  | 97       |
| 16) Acetophenone              | 8.745  | 105  | 297611   | 39.832 | ng/u1  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.734  | 70   | 136976   | 39.891 | ng/u1  | 97       |
| 18) 4-Methylphenol            | 8.704  | 108  | 189599   | 44.300 | ng/u1  | 97       |
| 19) Hexachloroethane          | 8.986  | 117  | 83516    | 37.538 | ng/u1  | 98       |
| 22) Nitrobenzene              | 9.128  | 77   | 210919   | 28.364 | ng/u1  | 98       |
| 23) Isophorone                | 9.657  | 82   | 412832   | 36.688 | ng/u1  | 99       |
| 25) 2-Nitrophenol             | 9.839  | 139  | 136897   | 42.329 | ng/u1  | 97       |
| 26) 2,4-Dimethylphenol        | 9.904  | 107  | 225932   | 32.162 | ng/u1  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 10.133 | 93   | 246583   | 42.977 | ng/u1  | 99       |
| 29) 2,4-Dichlorophenol        | 10.375 | 162  | 268289   | 39.091 | ng/u1  | 100      |
| 30) Naphthalene               | 10.769 | 128  | 664361   | 41.080 | ng/u1  | 99       |
| 32) 4-Chloroaniline           | 10.892 | 127  | 296288   | 42.076 | ng/u1  | 99       |
| 33) Hexachlorobutadiene       | 11.039 | 225  | 136668   | 19.161 | ng/u1  | 98       |
| 34) Caprolactam               | 11.698 | 113  | 60451m   | 44.769 | ng/u1  |          |
| 35) 4-Chloro-3-methylphenol   | 12.022 | 107  | 207379   | 38.252 | ng/u1  | 99       |
| 36) 2-Methylnaphthalene       | 12.374 | 142  | 524914   | 43.639 | ng/u1  | 99       |

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 Data File : BM038697.D  
 Acq On : 06 Feb 2023 17:54  
 Operator : CG/JU  
 Sample : SST04030  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST040036

Quant Time: Feb 07 00:53:10 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM020623.M  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 37) 1-Methylnaphthalene       | 12.598 | 142  | 539591   | 43.456 | ng/u1 | 97       |
| 39) 1,2,4,5-Tetrachloroben... | 12.745 | 216  | 253078   | 19.472 | ng/u1 | 97       |
| 40) Hexachlorocyclopentadiene | 12.710 | 237  | 138924   | 14.979 | ng/u1 | 92       |
| 41) 2,4,6-Trichlorophenol     | 12.992 | 196  | 206813   | 26.136 | ng/u1 | 95       |
| 42) 2,4,5-Trichlorophenol     | 13.069 | 196  | 224649   | 26.584 | ng/u1 | 97       |
| 43) 1,1'-Biphenyl             | 13.386 | 154  | 805766   | 39.690 | ng/u1 | 99       |
| 44) 2-Chloronaphthalene       | 13.427 | 162  | 650696   | 38.582 | ng/u1 | 99       |
| 45) 2-Nitroaniline            | 13.651 | 65   | 128592   | 28.506 | ng/u1 | 94       |
| 47) Dimethylphthalate         | 14.016 | 163  | 849505   | 39.584 | ng/u1 | 100      |
| 48) 2,6-Dinitrotoluene        | 14.145 | 165  | 173233   | 43.585 | ng/u1 | 99       |
| 50) Acenaphthylene            | 14.274 | 152  | 1002939  | 43.213 | ng/u1 | 98       |
| 51) 3-Nitroaniline            | 14.474 | 138  | 143268   | 50.077 | ng/u1 | 95       |
| 52) Acenaphthene              | 14.616 | 153  | 711225   | 41.170 | ng/u1 | 99       |
| 53) 2,4-Dinitrophenol         | 14.680 | 184  | 104557   | 32.723 | ng/u1 | 97       |
| 55) 4-Nitrophenol             | 14.798 | 109  | 121034   | 23.911 | ng/u1 | 97       |
| 56) Dibenzofuran              | 14.951 | 168  | 1005689  | 39.141 | ng/u1 | 100      |
| 57) 2,4-Dinitrotoluene        | 14.927 | 165  | 253743   | 43.313 | ng/u1 | 100      |
| 58) 2,3,4,6-Tetrachlorophenol | 15.180 | 232  | 169574   | 22.919 | ng/u1 | 98       |
| 59) Diethylphthalate          | 15.368 | 149  | 855999   | 42.547 | ng/u1 | 99       |
| 61) Fluorene                  | 15.598 | 166  | 832039   | 38.451 | ng/u1 | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.586 | 204  | 356514   | 23.415 | ng/u1 | 98       |
| 63) 4-Nitroaniline            | 15.639 | 138  | 119728   | 50.760 | ng/u1 | 95       |
| 66) 4,6-Dinitro-2-methylph... | 15.692 | 198  | 135218   | 29.485 | ng/u1 | 99       |
| 67) N-Nitrosodiphenylamine    | 15.810 | 169  | 694199   | 43.374 | ng/u1 | 99       |
| 68) 4-Bromophenyl-phenylether | 16.480 | 248  | 216802   | 27.104 | ng/u1 | 98       |
| 69) Hexachlorobenzene         | 16.592 | 284  | 277076   | 29.674 | ng/u1 | 97       |
| 70) Atrazine                  | 16.762 | 200  | 243426   | 30.312 | ng/u1 | 99       |
| 71) Pentachlorophenol         | 16.945 | 266  | 162612   | 27.225 | ng/u1 | 98       |
| 72) Phenanthrene              | 17.339 | 178  | 1296510  | 43.702 | ng/u1 | 99       |
| 74) Anthracene                | 17.427 | 178  | 1334925  | 42.970 | ng/u1 | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.351 | 216  | 253164   | 23.202 | ng/uL | 98       |
| 76) Pentachlorobenzene        | 14.863 | 250  | 281281   | 24.774 | ng/uL | 99       |
| 77) Carbazole                 | 17.704 | 167  | 1218097  | 45.172 | ng/u1 | 99       |
| 78) Di-n-butylphthalate       | 18.251 | 149  | 1610647  | 61.736 | ng/u1 | 100      |
| 80) Fluoranthene              | 19.351 | 202  | 1412009  | 50.017 | ng/u1 | 99       |
| 82) Pyrene                    | 19.715 | 202  | 1470883  | 49.096 | ng/u1 | 98       |
| 83) Butylbenzylphthalate      | 20.603 | 149  | 760124   | 87.962 | ng/u1 | 94       |
| 84) 3,3'-Dichlorobenzidine    | 21.392 | 252  | 457536   | 39.402 | ng/u1 | 100      |
| 85) Benzo(a)anthracene        | 21.456 | 228  | 1355201  | 40.474 | ng/u1 | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.368 | 149  | 1138858  | 92.619 | ng/u1 | 100      |
| 87) Chrysene                  | 21.509 | 228  | 1258681  | 40.613 | ng/u1 | 100      |
| 89) Di-n-octyl phthalate      | 22.286 | 149  | 1931358  | 92.770 | ng/u1 | 100      |
| 90) Benzo(b)fluoranthene      | 23.127 | 252  | 1353777  | 44.545 | ng/u1 | 99       |
| 91) Benzo(k)fluoranthene      | 23.174 | 252  | 1321122  | 43.611 | ng/u1 | 98       |
| 93) Benzo(a)pyrene            | 23.750 | 252  | 1118155  | 39.729 | ng/u1 | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.338 | 276  | 1372806  | 30.858 | ng/u1 | 99       |
| 95) Dibenzo(a,h)anthracene    | 26.344 | 278  | 1169479  | 30.974 | ng/u1 | 96       |
| 96) Benzo(g,h,i)perylene      | 27.097 | 276  | 1052109  | 29.137 | ng/u1 | 99       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM020623\  
 Data File : BM038697.D  
 Acq On : 06 Feb 2023 17:54  
 Operator : CG/JU  
 Sample : SSTD04030  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

## Instrument :

BNA\_M

## ClientSampleId :

SSTD040036

## Manual Integrations

## APPROVED

Reviewed By :Jagrut Upadhyay 02/07/2023  
 Supervised By :mohammad ahmed 02/07/2023

Quant Time: Feb 07 00:53:10 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM020623.M

Quant Title : SVOA CALIBRATION

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