

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061722\  
 Data File : BN020239.D  
 Acq On : 17 Jun 2022 14:56  
 Operator : CG/JU  
 Sample : N3313-04MSD  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 YB7L2MSD

Quant Time: Jun 17 22:56:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN060622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 22:53:31 2022  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 06/18/2022  
 Supervised By :mohammad ahmed 06/22/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.793  | 152  | 132417   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.581 | 136  | 656550   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.422 | 164  | 428817   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.163 | 188  | 865079   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.357 | 240  | 758196   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.668 | 264  | 677893   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.252  | 96   | 15417    | 5.043  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.664  | 84   | 90178    | 10.792 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.987  | 99   | 92052    | 8.401  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.122  | 67   | 239743   | 35.020 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.334  | 132  | 253444   | 28.468 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.516  | 113  | 185498   | 19.568 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.946  | 128  | 173736   | 34.783 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.663  | 143  | 185907   | 35.070 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.222 | 165  | 317801   | 34.342 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.716 | 131  | 460131   | 30.704 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.834 | 166  | 1153627  | 36.973 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.116 | 160  | 1315159  | 36.346 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.669 | 143  | 48122    | 7.697  | ng/u1  | 0.01     |
| 60) Fluorene-d10              | 15.416 | 176  | 960803   | 37.213 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.539 | 200  | 202839   | 40.478 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.263 | 188  | 1458952  | 38.643 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.563 | 212  | 1630836  | 39.677 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.527 | 264  | 1325109  | 39.952 | ng/u1  | 0.01     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.287  | 88   | 21456    | 6.648  | ng/uL# | 90       |
| 5) Pyridine                   | 3.681  | 79   | 99124    | 11.315 | ng/u1  | 90       |
| 6) Benzaldehyde               | 6.928  | 77   | 247862m  | 48.009 | ng/u1  |          |
| 8) Phenol                     | 7.011  | 94   | 105685   | 9.058  | ng/u1  | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.216  | 93   | 320735   | 35.122 | ng/u1  | 99       |
| 12) 2-Chlorophenol            | 7.363  | 128  | 265904   | 28.574 | ng/u1  | 98       |
| 13) 2-Methylphenol            | 8.252  | 108  | 206794   | 23.055 | ng/u1  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.322  | 45   | 507574   | 35.445 | ng/u1  | 97       |
| 16) Acetophenone              | 8.610  | 105  | 539484   | 37.421 | ng/u1  | 97       |
| 17) N-Nitroso-di-n-propyla... | 8.599  | 70   | 276771   | 37.582 | ng/u1  | 96       |
| 18) 4-Methylphenol            | 8.581  | 108  | 201408   | 20.345 | ng/u1  | 98       |
| 19) Hexachloroethane          | 8.869  | 117  | 126896   | 33.889 | ng/u1  | 86       |
| 22) Nitrobenzene              | 8.987  | 77   | 403646   | 34.853 | ng/u1  | 97       |
| 23) Isophorone                | 9.510  | 82   | 801064   | 35.032 | ng/u1  | 98       |
| 25) 2-Nitrophenol             | 9.699  | 139  | 202099   | 34.502 | ng/u1  | 97       |
| 26) 2,4-Dimethylphenol        | 9.775  | 107  | 305741   | 25.195 | ng/u1  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 9.999  | 93   | 485759   | 34.713 | ng/u1  | 97       |
| 29) 2,4-Dichlorophenol        | 10.246 | 162  | 322050   | 33.643 | ng/u1  | 98       |
| 30) Naphthalene               | 10.634 | 128  | 1155350  | 33.954 | ng/u1  | 99       |
| 32) 4-Chloroaniline           | 10.740 | 127  | 452412   | 30.182 | ng/u1  | 100      |
| 33) Hexachlorobutadiene       | 10.928 | 225  | 178281   | 32.439 | ng/u1  | 97       |
| 34) Caprolactam               | 11.499 | 113  | 17982    | 4.810  | ng/u1  | 99       |
| 35) 4-Chloro-3-methylphenol   | 11.887 | 107  | 367586   | 32.521 | ng/u1  | 99       |
| 36) 2-Methylnaphthalene       | 12.246 | 142  | 816723   | 35.503 | ng/u1  | 98       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061722\  
 Data File : BN020239.D  
 Acq On : 17 Jun 2022 14:56  
 Operator : CG/JU  
 Sample : N3313-04MSD  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

## Instrument :

BNA\_N

## ClientSampleId :

YB7L2MSD

## Manual Integrations

## APPROVED

Reviewed By :Jagrut Upadhyay 06/18/2022  
 Supervised By :mohammad ahmed 06/22/2022

Quant Time: Jun 17 22:56:53 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN060622.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Jun 17 22:53:31 2022

Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.463 | 142  | 846454   | 36.057 | ng/ul  | 94       |
| 39) 1,2,4,5-Tetrachloroben... | 12.616 | 216  | 376167   | 34.395 | ng/ul  | 98       |
| 40) Hexachlorocyclopentadiene | 12.598 | 237  | 159043   | 23.906 | ng/ul  | 95       |
| 41) 2,4,6-Trichlorophenol     | 12.863 | 196  | 280205   | 37.157 | ng/ul  | 93       |
| 42) 2,4,5-Trichlorophenol     | 12.945 | 196  | 304811   | 36.496 | ng/ul  | 98       |
| 43) 1,1'-Biphenyl             | 13.257 | 154  | 1120199  | 35.279 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.298 | 162  | 847871   | 35.198 | ng/ul  | 96       |
| 45) 2-Nitroaniline            | 13.504 | 65   | 299287   | 39.937 | ng/ul  | 93       |
| 47) Dimethylphthalate         | 13.881 | 163  | 1134351  | 36.114 | ng/ul  | 98       |
| 48) 2,6-Dinitrotoluene        | 13.998 | 165  | 248344   | 40.439 | ng/ul  | 97       |
| 50) Acenaphthylene            | 14.145 | 152  | 1436675  | 35.862 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.328 | 138  | 244346   | 37.635 | ng/ul  | 98       |
| 52) Acenaphthene              | 14.487 | 153  | 926568   | 35.409 | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol         | 14.534 | 184  | 137295   | 34.918 | ng/ul  | 94       |
| 55) 4-Nitrophenol             | 14.681 | 109  | 42093    | 7.887  | ng/ul  | 93       |
| 56) Dibenzofuran              | 14.822 | 168  | 1342149  | 36.406 | ng/ul  | 97       |
| 57) 2,4-Dinitrotoluene        | 14.787 | 165  | 356596   | 39.216 | ng/ul  | 93       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.057 | 232  | 242406   | 36.974 | ng/ul  | 92       |
| 59) Diethylphthalate          | 15.245 | 149  | 1209943  | 37.415 | ng/ul  | 97       |
| 61) Fluorene                  | 15.469 | 166  | 1051073  | 36.639 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.463 | 204  | 477007   | 36.107 | ng/ul  | 91       |
| 63) 4-Nitroaniline            | 15.492 | 138  | 252841   | 41.676 | ng/ul  | 98       |
| 66) 4,6-Dinitro-2-methylph... | 15.551 | 198  | 202032   | 39.766 | ng/ul  | 94       |
| 67) N-Nitrosodiphenylamine    | 15.681 | 169  | 915688   | 36.922 | ng/ul  | 97       |
| 68) 4-Bromophenyl-phenylether | 16.357 | 248  | 294943   | 39.151 | ng/ul# | 85       |
| 69) Hexachlorobenzene         | 16.481 | 284  | 324980   | 37.527 | ng/ul# | 92       |
| 70) Atrazine                  | 16.634 | 200  | 341395   | 39.453 | ng/ul  | 96       |
| 71) Pentachlorophenol         | 16.822 | 266  | 196051   | 36.349 | ng/ul  | 97       |
| 72) Phenanthrene              | 17.204 | 178  | 1734920  | 38.002 | ng/ul  | 99       |
| 74) Anthracene                | 17.298 | 178  | 1735074  | 37.797 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.222 | 216  | 413641   | 35.707 | ng/uL  | 98       |
| 76) Pentachlorobenzene        | 14.745 | 250  | 382658   | 37.375 | ng/uL  | 96       |
| 77) Carbazole                 | 17.569 | 167  | 1692121  | 40.518 | ng/ul  | 98       |
| 78) Di-n-butylphthalate       | 18.139 | 149  | 2136880  | 41.264 | ng/ul  | 100      |
| 80) Fluoranthene              | 19.222 | 202  | 2006828  | 40.425 | ng/ul  | 96       |
| 82) Pyrene                    | 19.586 | 202  | 1973314  | 38.745 | ng/ul# | 92       |
| 83) Butylbenzylphthalate      | 20.492 | 149  | 1005497  | 43.868 | ng/ul  | 97       |
| 84) 3,3'-Dichlorobenzidine    | 21.274 | 252  | 495373   | 34.208 | ng/ul# | 94       |
| 85) Benzo(a)anthracene        | 21.345 | 228  | 1824832  | 38.115 | ng/ul  | 97       |
| 86) Bis(2-ethylhexyl)phtha... | 21.280 | 149  | 1415734  | 43.777 | ng/ul  | 99       |
| 87) Chrysene                  | 21.392 | 228  | 1774582  | 37.886 | ng/ul  | 100      |
| 89) Di-n-octyl phthalate      | 22.180 | 149  | 2554838  | 46.305 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 22.974 | 252  | 1686846  | 39.648 | ng/ul# | 96       |
| 91) Benzo(k)fluoranthene      | 23.016 | 252  | 1695031  | 41.852 | ng/ul  | 96       |
| 93) Benzo(a)pyrene            | 23.574 | 252  | 1596690  | 38.770 | ng/ul  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.045 | 276  | 1540953  | 36.367 | ng/ul# | 94       |
| 95) Dibenzo(a,h)anthracene    | 26.062 | 278  | 1334296  | 36.868 | ng/ul  | 96       |
| 96) Benzo(g,h,i)perylene      | 26.774 | 276  | 1327037  | 37.087 | ng/ul# | 90       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061722\  
 Data File : BN020239.D  
 Acq On : 17 Jun 2022 14:56  
 Operator : CG/JU  
 Sample : N3313-04MSD  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 YB7L2MSD

### Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 06/18/2022  
 Supervised By :mohammad ahmed 06/22/2022

Quant Time: Jun 17 22:56:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN060622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 22:53:31 2022  
 Response via : Initial Calibration

