

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM120922\  
 Data File : BM037917.D  
 Acq On : 09 Dec 2022 15:16  
 Operator : CG/JU  
 Sample : N5896-05 10X  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBXM3

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2022  
 Supervised By :mohammad ahmed 12/10/2022

Quant Time: Dec 09 21:48:41 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 09 21:34:57 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.081  | 152  | 267835   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.904 | 136  | 1101223  | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.710 | 164  | 635347   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.451 | 188  | 1315402  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.621 | 240  | 969847   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.109 | 264  | 1061611  | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.422  | 96   | 1689m    | 0.286  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.887  | 84   | 13812m   | 0.789  | ng/u1  | 0.03     |
| 7) Phenol-d5                  | 7.234  | 99   | 37710    | 1.733  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.404  | 67   | 23595    | 1.780  | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.610  | 132  | 32479    | 1.819  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.792  | 113  | 28396    | 1.671  | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.251  | 128  | 15780    | 1.810  | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.981  | 143  | 14725    | 1.612  | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.522 | 165  | 27142    | 1.564  | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.039 | 131  | 28127    | 1.174  | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.116 | 166  | 84003    | 1.855  | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.404 | 160  | 96275    | 1.739  | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.910 | 143  | 7114m    | 0.922  | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.698 | 176  | 76094    | 1.886  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.816 | 200  | 6916     | 0.983  | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.551 | 188  | 130698   | 2.129  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.833 | 212  | 129469   | 2.222  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.939 | 264  | 106219   | 2.003  | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 30) Naphthalene               | 10.951 | 128  | 85793    | 1.444  | ng/u1  | 99       |
| 50) Acenaphthylene            | 14.439 | 152  | 121779   | 2.001  | ng/u1  | 99       |
| 52) Acenaphthene              | 14.774 | 153  | 48961    | 1.163  | ng/u1  | 96       |
| 61) Fluorene                  | 15.751 | 166  | 104749   | 2.234  | ng/u1  | 98       |
| 72) Phenanthrene              | 17.492 | 178  | 696706   | 9.699  | ng/u1  | 100      |
| 74) Anthracene                | 17.586 | 178  | 2918986  | 39.985 | ng/u1  | 99       |
| 80) Fluoranthene              | 19.498 | 202  | 2150370  | 31.275 | ng/u1  | 100      |
| 82) Pyrene                    | 19.868 | 202  | 5268359  | 73.624 | ng/u1  | 99       |
| 85) Benzo(a)anthracene        | 21.603 | 228  | 630869   | 9.662  | ng/u1  | 97       |
| 87) Chrysene                  | 21.662 | 228  | 1131965m | 18.478 | ng/u1  |          |
| 90) Benzo(b)fluoranthene      | 23.350 | 252  | 1903852  | 28.927 | ng/u1  | 98       |
| 91) Benzo(k)fluoranthene      | 23.392 | 252  | 338312m  | 5.304  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 23.997 | 252  | 771973   | 13.338 | ng/u1  | 97       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.697 | 276  | 439179   | 5.892  | ng/u1# | 93       |
| 95) Dibenzo(a,h)anthracene    | 26.697 | 278  | 108596   | 1.728  | ng/u1# | 85       |
| 96) Benzo(g,h,i)perylene      | 27.497 | 276  | 329169   | 5.531  | ng/u1  | 98       |
| -----                         |        |      |          |        |        |          |

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

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