

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM120922\  
 Data File : BM037910.D  
 Acq On : 09 Dec 2022 10:59  
 Operator : CG/JU  
 Sample : N5832-16DL 20X  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBXL1DL

Manual Integrations  
 APPROVED

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

Quant Time: Dec 09 21:46:59 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  | 251734   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.904 | 136  | 992996   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.710 | 164  | 536922   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.451 | 188  | 1039727  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.621 | 240  | 891818   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.103 | 264  | 999996   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.428  | 96   | 1328m    | 0.240  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 7.234  | 99   | 17425    | 0.852  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.404  | 67   | 12089    | 0.971  | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.610  | 132  | 15214    | 0.907  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.792  | 113  | 13325    | 0.834  | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.257  | 128  | 7315     | 0.930  | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.980  | 143  | 6230     | 0.756  | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.522 | 165  | 12225    | 0.781  | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.051 | 131  | 9783     | 0.453  | ng/u1  | 0.01     |
| 46) Dimethylphthalate-d6      | 14.115 | 166  | 37239    | 0.973  | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.404 | 160  | 43523    | 0.930  | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000  | ng/u1  |          |
| 60) Fluorene-d10              | 15.698 | 176  | 36222    | 1.062  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.821 | 200  | 2043     | 0.368  | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.551 | 188  | 63627    | 1.311  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.833 | 212  | 61253    | 1.143  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.938 | 264  | 61334    | 1.228  | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 30) Naphthalene               | 10.951 | 128  | 71108    | 1.327  | ng/u1  | 99       |
| 36) 2-Methylnaphthalene       | 12.551 | 142  | 43901    | 1.263  | ng/u1  | 98       |
| 50) Acenaphthylene            | 14.439 | 152  | 57312    | 1.114  | ng/u1  | 98       |
| 52) Acenaphthene              | 14.774 | 153  | 64168    | 1.803  | ng/u1  | 98       |
| 61) Fluorene                  | 15.757 | 166  | 199403   | 5.033  | ng/u1  | 97       |
| 72) Phenanthrene              | 17.492 | 178  | 527267   | 9.286  | ng/u1  | 99       |
| 74) Anthracene                | 17.586 | 178  | 3199182  | 55.443 | ng/u1  | 98       |
| 80) Fluoranthene              | 19.497 | 202  | 413211   | 6.536  | ng/u1  | 98       |
| 82) Pyrene                    | 19.862 | 202  | 574185   | 8.726  | ng/u1  | 97       |
| 85) Benzo(a)anthracene        | 21.603 | 228  | 134240   | 2.236  | ng/u1  | 96       |
| 87) Chrysene                  | 21.656 | 228  | 281711m  | 5.001  | ng/u1  |          |
| 90) Benzo(b)fluoranthene      | 23.344 | 252  | 608566m  | 9.816  | ng/u1  |          |
| 91) Benzo(k)fluoranthene      | 23.385 | 252  | 118905m  | 1.979  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 23.991 | 252  | 243398   | 4.465  | ng/u1  | 95       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.691 | 276  | 175686   | 2.502  | ng/u1  | 93       |
| 96) Benzo(g,h,i)perylene      | 27.497 | 276  | 129321   | 2.307  | ng/u1  | 94       |
| -----                         |        |      |          |        |        |          |

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

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