

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM121823\  
 Data File : BM043395.D  
 Acq On : 18 Dec 2023 16:19  
 Operator : MA/JU  
 Sample : 05626-16DL 5X  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DCLG2DL

Quant Time: Dec 19 01:24:18 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 15 22:09:25 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 12/19/2023  
 Supervised By :mohammad ahmed 12/20/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.993  | 152  | 517705   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.804 | 136  | 2243306  | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.621 | 164  | 1155117  | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.368 | 188  | 1938558  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.545 | 240  | 1333664  | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.974 | 264  | 1487896  | 20.000 | ng/u1  | 0.02     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.387  | 96   | 11073    | 0.695  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 50431    | 1.100  | ng/u1  | 0.02     |
| 7) Phenol-d5                  | 7.151  | 99   | 222667   | 3.781  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.328  | 67   | 146936   | 3.940  | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.522  | 132  | 171473   | 3.813  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.704  | 113  | 176896   | 3.871  | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.163  | 128  | 83210    | 3.834  | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.887  | 143  | 82205    | 3.666  | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.422 | 165  | 149171   | 3.768  | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.945 | 131  | 187505   | 3.219  | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.039 | 166  | 428365   | 4.120  | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.316 | 160  | 497229   | 4.031  | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.821 | 143  | 48280    | 2.453  | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.616 | 176  | 345179   | 4.014  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.733 | 200  | 30820    | 2.404  | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.462 | 188  | 449454   | 4.134  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.751 | 212  | 427154   | 4.057  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.809 | 264  | 384457   | 4.187  | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 30) Naphthalene               | 10.851 | 128  | 3675774  | 25.710 | ng/u1  | 100      |
| 36) 2-Methylnaphthalene       | 12.457 | 142  | 263431   | 2.707  | ng/u1  | 100      |
| 37) 1-Methylnaphthalene       | 12.675 | 142  | 183730   | 1.872  | ng/u1  | 99       |
| 74) Anthracene                | 17.498 | 178  | 144840   | 1.092  | ng/u1  | 98       |
| 80) Fluoranthene              | 19.421 | 202  | 346909   | 2.661  | ng/u1  | 99       |
| 82) Pyrene                    | 19.780 | 202  | 416891   | 3.110  | ng/u1  | 98       |
| 85) Benzo(a)anthracene        | 21.527 | 228  | 193978   | 1.739  | ng/u1  | 92       |
| 87) Chrysene                  | 21.556 | 228  | 377487m  | 3.763  | ng/u1  |          |
| 90) Benzo(b)fluoranthene      | 23.233 | 252  | 846052   | 7.420  | ng/u1  | 97       |
| 91) Benzo(k)fluoranthene      | 23.274 | 252  | 252474m  | 2.270  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 23.862 | 252  | 320691   | 3.083  | ng/u1  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.491 | 276  | 265801   | 2.140  | ng/u1  | 98       |
| -----                         |        |      |          |        |        |          |

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

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