

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM121523\  
 Data File : BM043377.D  
 Acq On : 16 Dec 2023 02:51  
 Operator : MA/JU  
 Sample : 05626-20  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DCLG6

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 12/16/2023  
 Supervised By :mohammad ahmed 12/16/2023

Quant Time: Dec 16 03:24:27 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

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.992  | 152  | 423392   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.804 | 136  | 1765261  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.621 | 164  | 954169   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.368 | 188  | 1826300  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.539 | 240  | 1475103  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.962 | 264  | 1742506  | 20.000 | ng/u1 | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.387  | 96   | 49330    | 3.787  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.810  | 84   | 309267   | 8.249  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.151  | 99   | 975068   | 20.245 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.334  | 67   | 619043   | 20.294 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.522  | 132  | 765041   | 20.804 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.704  | 113  | 700659   | 18.747 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.169  | 128  | 367170   | 21.498 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.886  | 143  | 384220   | 21.773 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.422 | 165  | 654079   | 20.996 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.945 | 131  | 785474   | 17.139 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.045 | 166  | 1847242  | 21.506 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.321 | 160  | 2256392  | 22.144 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.821 | 143  | 306069   | 18.827 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.615 | 176  | 1605309  | 22.601 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.733 | 200  | 176212   | 14.589 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.468 | 188  | 2282076  | 22.281 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.750 | 212  | 2529742  | 21.722 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.809 | 264  | 2463324  | 22.908 | ng/u1 | 0.00     |
|                               |        |      |          |        |       |          |
| Target Compounds              |        |      |          |        |       | Qvalue   |
| 30) Naphthalene               | 10.851 | 128  | 153459   | 1.364  | ng/u1 | 97       |
| 52) Acenaphthene              | 14.686 | 153  | 276418   | 3.584  | ng/u1 | 100      |
| 61) Fluorene                  | 15.668 | 166  | 334100   | 3.927  | ng/u1 | 100      |
| 72) Phenanthrene              | 17.409 | 178  | 2007895  | 16.222 | ng/u1 | 100      |
| 74) Anthracene                | 17.498 | 178  | 895347   | 7.167  | ng/u1 | 99       |
| 80) Fluoranthene              | 19.421 | 202  | 3037564  | 21.062 | ng/u1 | 97       |
| 82) Pyrene                    | 19.780 | 202  | 2380139  | 16.054 | ng/u1 | 100      |
| 85) Benzo(a)anthracene        | 21.521 | 228  | 645892   | 5.236  | ng/u1 | 97       |
| 87) Chrysene                  | 21.574 | 228  | 770092   | 6.940  | ng/u1 | 99       |
| 90) Benzo(b)fluoranthene      | 23.221 | 252  | 1049238  | 7.858  | ng/u1 | 98       |
| 91) Benzo(k)fluoranthene      | 23.268 | 252  | 295095m  | 2.266  | ng/u1 |          |
| 93) Benzo(a)pyrene            | 23.850 | 252  | 363592   | 2.984  | ng/u1 | 96       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.479 | 276  | 278674   | 1.916  | ng/u1 | 98       |
| -----                         |        |      |          |        |       |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM121523\  
Data File : BM043377.D  
Acq On : 16 Dec 2023 02:51  
Operator : MA/JU  
Sample : 05626-20  
Misc :  
ALS Vial : 26 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
DCLG6

Quant Time: Dec 16 03:24:27 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 12/16/2023  
Supervised By :mohammad ahmed 12/16/2023

