

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062923\  
 Data File : BP015821.D  
 Acq On : 29 Jun 2023 21:28  
 Operator : MA/JU  
 Sample : 03143-09  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCFU6

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 06/30/2023  
 Supervised By :mohammad ahmed 06/30/2023

Quant Time: Jun 30 00:07:31 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062623.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 29 02:30:23 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.058  | 152  | 263217   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.887 | 136  | 1141385  | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.704 | 164  | 679010   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.463 | 188  | 1420820  | 20.000 | ng/u1  | # 0.00   |
| 79) Chrysene-d12              | 21.551 | 240  | 1036295  | 20.000 | ng/u1  | # 0.00   |
| 88) Perylene-d12              | 24.098 | 264  | 1005612  | 20.000 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.381  | 96   | 21919    | 2.996  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 7.234  | 99   | 420715   | 18.681 | ng/u1  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.381  | 67   | 287083   | 20.604 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.587  | 132  | 342440   | 20.143 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.793  | 113  | 265522   | 14.924 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.246  | 128  | 166644   | 19.104 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.975  | 143  | 180712   | 17.750 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.546 | 165  | 303858   | 16.749 | ng/u1  | 0.03     |
| 31) 4-Chloroaniline-d4        | 11.069 | 131  | 169517   | 7.274  | ng/u1  | 0.04     |
| 46) Dimethylphthalate-d6      | 14.116 | 166  | 1078219  | 20.544 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.398 | 160  | 1233636  | 20.910 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.993 | 143  | 91763m   | 13.249 | ng/u1  | 0.04     |
| 60) Fluorene-d10              | 15.704 | 176  | 919643   | 21.281 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.869 | 200  | 87601m   | 10.025 | ng/u1  | 0.01     |
| 73) Anthracene-d10            | 17.569 | 188  | 1314442  | 20.253 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.792 | 212  | 1574412  | 23.267 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.927 | 264  | 1105678  | 21.796 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 80) Fluoranthene              | 19.469 | 202  | 177684   | 2.113  | ng/u1# | 91       |
| 82) Pyrene                    | 19.822 | 202  | 151221   | 1.763  | ng/u1# | 87       |
| 85) Benzo(a)anthracene        | 21.539 | 228  | 81825    | 1.122  | ng/u1  | 96       |
| 87) Chrysene                  | 21.592 | 228  | 99367    | 1.437  | ng/u1  | 95       |
| 90) Benzo(b)fluoranthene      | 23.316 | 252  | 142438   | 2.204  | ng/u1# | 79       |
| 93) Benzo(a)pyrene            | 23.986 | 252  | 86547    | 1.533  | ng/u1# | 89       |
| 96) Benzo(g,h,i)perylene      | 27.645 | 276  | 69333    | 1.100  | ng/u1# | 88       |
| -----                         |        |      |          |        |        |          |

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

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