

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG082124\  
 Data File : BG062526.D  
 Acq On : 22 Aug 2024 3:32  
 Operator : MA/JU  
 Sample : P3626-03  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DCYC7

Quant Time: Aug 22 04:14:40 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG081424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Aug 20 23:01:24 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.939  | 152  | 83368    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.730 | 136  | 389435   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.560 | 164  | 290308   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.297 | 188  | 619270   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.533 | 240  | 597709   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 24.611 | 264  | 698683   | 20.000 | ng/u1 | -0.01    |

|                               |        |     |        |        |       |      |
|-------------------------------|--------|-----|--------|--------|-------|------|
| System Monitoring Compounds   |        |     |        |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.404  | 96  | 8446   | 3.846  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.810  | 84  | 140496 | 20.997 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.099  | 99  | 203717 | 24.563 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.264  | 67  | 117939 | 22.576 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.470  | 132 | 146723 | 25.899 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.639  | 113 | 165209 | 23.765 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.091  | 128 | 74551  | 29.851 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.813  | 143 | 82709  | 33.292 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.348 | 165 | 167522 | 26.474 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.859 | 131 | 229288 | 22.849 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.961 | 166 | 535092 | 23.696 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.249 | 160 | 613599 | 24.541 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.742 | 143 | 81716  | 22.684 | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.547 | 176 | 500097 | 24.554 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.658 | 200 | 70352  | 25.003 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.397 | 188 | 750423 | 25.225 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.682 | 212 | 921080 | 26.313 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 24.405 | 264 | 954921 | 25.754 | ng/u1 | 0.00 |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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