

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110722\  
 Data File : BP012538.D  
 Acq On : 07 Nov 2022 19:07  
 Operator : CG/JU  
 Sample : N5353-13  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DBWW9

Quant Time: Nov 07 21:42:20 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 04 02:53:36 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.775  | 152  | 304397   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.575 | 136  | 1458888  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.428 | 164  | 996357   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.192 | 188  | 2242936  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.286 | 240  | 2081549  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.669 | 264  | 1914966  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.228  | 96   | 27550    | 3.644  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.658  | 84   | 153827   | 7.083  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.958  | 99   | 146866   | 5.379  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.117  | 67   | 507920   | 31.164 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.311  | 132  | 468273   | 23.487 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.499  | 113  | 321757   | 14.901 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.946  | 128  | 354753   | 37.498 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.663  | 143  | 365647   | 37.237 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.205 | 165  | 629292   | 29.571 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.728 | 131  | 789589   | 24.919 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.846 | 166  | 2413046  | 35.379 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.122 | 160  | 2625942  | 33.690 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.681 | 143  | 37790    | 2.927  | ng/u1 | 0.02     |
| 60) Fluorene-d10              | 15.428 | 176  | 2115827  | 36.232 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.563 | 200  | 324016   | 27.747 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.292 | 188  | 3529510  | 36.521 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.534 | 212  | 4135845  | 39.575 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.516 | 264  | 3474769  | 37.136 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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