

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP041122\  
 Data File : BP009788.D  
 Acq On : 12 Apr 2022 18:08  
 Operator : CG/JU  
 Sample : N2342-03  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0J67

Quant Time: Apr 13 02:24:12 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP040722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Apr 13 02:15:22 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.963  | 152  | 98039    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.775 | 136  | 423541   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.598 | 164  | 296617   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.345 | 188  | 665816   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.427 | 240  | 713186   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.892 | 264  | 692642   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.387  | 96   | 11568    | 5.266  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.805  | 84   | 62777    | 10.365 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.116  | 99   | 62533    | 7.265  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.287  | 67   | 188430   | 36.773 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.493  | 132  | 177440   | 27.330 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.663  | 113  | 122658   | 17.071 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.122  | 128  | 116921   | 38.281 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.851  | 143  | 122419   | 36.559 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.387 | 165  | 213398   | 30.455 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.904 | 131  | 302787   | 30.339 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.004 | 166  | 897407   | 38.516 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.286 | 160  | 973748   | 35.276 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.769 | 143  | 29726    | 7.035  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.586 | 176  | 779587   | 39.774 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.698 | 200  | 130744   | 33.147 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.445 | 188  | 1310795  | 41.118 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.674 | 212  | 1687712  | 43.329 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.733 | 264  | 1596887  | 44.271 | ng/u1 | 0.00     |

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

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

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