

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080922\  
 Data File : BP011328.D  
 Acq On : 09 Aug 2022 14:20  
 Operator : CG/JU  
 Sample : N4026-21  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Aug 09 23:11:13 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 08 18:47:19 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.593  | 152  | 121603   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.381 | 136  | 512765   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.263 | 164  | 289392   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.033 | 188  | 578632   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.163 | 240  | 532351   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.474 | 264  | 527786   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.164  | 96   | 79       | 0.020  | ng/uL | 0.08     |
| 4) Pyridine-d5                | 3.517  | 84   | 45187    | 4.516  | ng/u1 | 0.02     |
| 7) Phenol-d5                  | 6.775  | 99   | 44274    | 4.114  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.940  | 67   | 188634   | 24.721 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.122  | 132  | 144879   | 17.158 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.310  | 113  | 86837    | 10.277 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.757  | 128  | 101610   | 29.117 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.475  | 143  | 81366    | 25.466 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.010 | 165  | 136822   | 21.205 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.534 | 131  | 203136   | 18.008 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.687 | 166  | 599299   | 29.846 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.951 | 160  | 688908   | 28.502 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.498 | 143  | 4594     | 1.253  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.269 | 176  | 494019   | 28.769 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.398 | 200  | 49921    | 18.574 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.133 | 188  | 806284   | 31.001 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.398 | 212  | 906013   | 30.891 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.327 | 264  | 818393   | 30.747 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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