

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121922\  
 Data File : BP013227.D  
 Acq On : 22 Dec 2022 02:35  
 Operator : CG/JU  
 Sample : N6032-15  
 Misc :  
 ALS Vial : 110 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0LC2

Quant Time: Dec 22 05:39:13 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 21 23:38:37 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.945  | 152  | 536842   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.763 | 136  | 2100516  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.592 | 164  | 1205810  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.345 | 188  | 2587566  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.433 | 240  | 2806008  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.909 | 264  | 3316463  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.334  | 96   | 50288    | 4.005  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.775  | 84   | 108003   | 3.028  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 7.104  | 99   | 304060   | 6.954  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 701261   | 26.585 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.475  | 132  | 576418   | 16.745 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.657  | 113  | 374603   | 10.455 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.116  | 128  | 443807   | 28.757 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.839  | 143  | 398817   | 22.954 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.381 | 165  | 662663   | 19.636 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.904 | 131  | 204551   | 4.293  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.998 | 166  | 2629305  | 28.372 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.286 | 160  | 3009329  | 29.552 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.786 | 143  | 64666    | 3.879  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.580 | 176  | 2419572  | 30.862 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.704 | 200  | 285182   | 17.126 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.445 | 188  | 3932197  | 33.684 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.674 | 212  | 5043996  | 31.317 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.750 | 264  | 5436397  | 32.887 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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