

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061323\  
 Data File : BP015499.D  
 Acq On : 13 Jun 2023 16:43  
 Operator : MA/JU  
 Sample : 03097-06  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EZEN3

Quant Time: Jun 14 00:56:07 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 09 23:33:38 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.093  | 152  | 177799   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.916 | 136  | 819535   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.734 | 164  | 550357   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.492 | 188  | 1265566  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.575 | 240  | 1201366  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.122 | 264  | 1256109  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.411  | 96   | 20059    | 4.178  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.858  | 84   | 75820    | 6.079  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 7.246  | 99   | 104137   | 6.358  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.411  | 67   | 326715   | 33.398 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.616  | 132  | 303079   | 25.520 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.805  | 113  | 216323   | 16.092 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.269  | 128  | 214902   | 33.400 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.999  | 143  | 235927   | 32.109 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.540 | 165  | 391978   | 29.244 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.057 | 131  | 502145   | 26.100 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.140 | 166  | 1605800  | 37.001 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.428 | 160  | 1726378  | 36.379 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.963 | 143  | 32949    | 4.240  | ng/u1 | 0.02     |
| 60) Fluorene-d10              | 15.728 | 176  | 1345468  | 36.764 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.851 | 200  | 262341   | 32.721 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.592 | 188  | 2255622  | 39.173 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.816 | 212  | 2702832  | 42.035 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.957 | 264  | 2524838  | 40.060 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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