

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP050923\  
 Data File : BP015036.D  
 Acq On : 10 May 2023 01:21  
 Operator : CG/JU  
 Sample : 02694-01  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 JREL2

Quant Time: May 10 02:54:40 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP050323.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri May 05 12:12:36 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.157  | 152  | 284327   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.993 | 136  | 1295613  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.798 | 164  | 698120   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.557 | 188  | 1525092  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.633 | 240  | 1250416  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 24.227 | 264  | 1067515  | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.452  | 96   | 34042    | 4.452  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.899  | 84   | 219314   | 10.493 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.299  | 99   | 252618   | 10.248 | ng/u1 | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.475  | 67   | 466267   | 30.602 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.681  | 132  | 508350   | 28.046 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.863  | 113  | 409154   | 21.538 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.334  | 128  | 313645   | 34.038 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.063 | 143  | 325103   | 34.359 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.604 | 165  | 602195   | 32.414 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 11.128 | 131  | 610452   | 22.717 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.204 | 166  | 1836977  | 37.179 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.492 | 160  | 2071841  | 34.963 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.987 | 143  | 81460    | 9.477  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.792 | 176  | 1535734  | 36.128 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.904 | 200  | 251073   | 29.245 | ng/u1 | -0.02    |
| 73) Anthracene-d10            | 17.657 | 188  | 2444247  | 35.670 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.875 | 212  | 2838249  | 33.655 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.057 | 264  | 2027229  | 38.975 | ng/u1 | -0.02    |

Target Compounds Qvalue

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

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