

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP041522\  
 Data File : BP009897.D  
 Acq On : 16 Apr 2022 12:21  
 Operator : CG/JU  
 Sample : N2421-18  
 Misc :  
 ALS Vial : 49 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0J80

Quant Time: Apr 18 01:12:43 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP040722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Apr 18 00:59:30 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.958  | 152  | 145971   | 20.000 | ng/ul | # 0.00   |
| 20) Naphthalene-d8            | 10.763 | 136  | 643835   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.587 | 164  | 460562   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.340 | 188  | 1038808  | 20.000 | ng/ul | # 0.00   |
| 79) Chrysene-d12              | 21.416 | 240  | 1050632  | 20.000 | ng/ul | # 0.00   |
| 88) Perylene-d12              | 23.880 | 264  | 998555   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.381  | 96   | 14755    | 4.511  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.799  | 84   | 77635    | 8.609  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.111  | 99   | 79206    | 6.181  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.281  | 67   | 221983   | 29.096 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.481  | 132  | 210042   | 21.728 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.658  | 113  | 156084   | 14.590 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.111  | 128  | 145526   | 31.344 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.834  | 143  | 149901   | 29.449 | ng/ul | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.375 | 165  | 259742   | 24.385 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.893 | 131  | 421666   | 27.794 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.993 | 166  | 1090880  | 30.153 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.281 | 160  | 1196760  | 27.922 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.763 | 143  | 35924    | 5.475  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 952534   | 31.298 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.687 | 200  | 156415   | 25.417 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.440 | 188  | 1671323  | 33.603 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.663 | 212  | 2055933  | 35.830 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.722 | 264  | 1860458  | 35.777 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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