

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081423\  
 Data File : BP016572.D  
 Acq On : 14 Aug 2023 21:49  
 Operator : MA/JU  
 Sample : PB154810BL  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK810

Quant Time: Aug 15 00:03:59 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 14 18:26:05 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 8.063  | 152  | 736619   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.893 | 136  | 3244291  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.710 | 164  | 1944432  | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.475 | 188  | 4130242  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.569 | 240  | 3489328  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12          | 24.157 | 264  | 3294375  | 20.000 | ng/ul | 0.00     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.399  | 96  | 69429   | 3.582  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.834  | 84  | 1045623 | 17.729 | ng/ul | 0.00 |
| 7) Phenol-d5                  | 7.199  | 99  | 1248801 | 18.389 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.393  | 67  | 787307  | 17.602 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.575  | 132 | 916339  | 18.786 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8         | 8.763  | 113 | 949269  | 17.580 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5           | 9.258  | 128 | 455338  | 19.947 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.975  | 143 | 475360  | 21.925 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.499 | 165 | 824268  | 19.318 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4        | 11.046 | 131 | 1281742 | 17.508 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6      | 14.122 | 166 | 2555884 | 18.173 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8         | 14.410 | 160 | 2974242 | 18.737 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.904 | 143 | 449763  | 15.698 | ng/ul | 0.00 |
| 60) Fluorene-d10              | 15.704 | 176 | 2185767 | 18.230 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.822 | 200 | 241895  | 12.311 | ng/ul | 0.00 |
| 73) Anthracene-d10            | 17.575 | 188 | 3380459 | 18.922 | ng/ul | 0.00 |
| 81) Pyrene-d10                | 19.804 | 212 | 3775343 | 20.649 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.986 | 264 | 3119237 | 19.757 | ng/ul | 0.00 |

Target Compounds Qvalue

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

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