

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122222\  
 Data File : BP013276.D  
 Acq On : 23 Dec 2022 13:09  
 Operator : CG/JU  
 Sample : N6015-21  
 Misc :  
 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DBYA4

Quant Time: Dec 24 00:24:42 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Dec 24 00:19:17 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.946  | 152  | 683107   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.757 | 136  | 2882365  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.592 | 164  | 1813679  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.351 | 188  | 3475155  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.439 | 240  | 2554324  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.927 | 264  | 2953118  | 20.000 | ng/u1 | 0.01     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.334  | 96  | 112566  | 7.046  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.758  | 84  | 1775746 | 39.127 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.110  | 99  | 2516791 | 45.233 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67  | 1484416 | 44.226 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.475  | 132 | 2002423 | 45.716 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.657  | 113 | 1992094 | 43.695 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.116  | 128 | 1030383 | 48.655 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.840  | 143 | 1215711 | 50.991 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.381 | 165 | 2206512 | 47.647 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.898 | 131 | 2608789 | 39.904 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.998 | 166 | 6535698 | 46.888 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.287 | 160 | 7456318 | 48.681 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.792 | 143 | 938619  | 37.435 | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.586 | 176 | 5485197 | 46.515 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.704 | 200 | 830921  | 37.155 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.451 | 188 | 7430606 | 47.395 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.680 | 212 | 7433623 | 50.701 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.768 | 264 | 7244608 | 49.218 | ng/u1 | 0.01 |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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