

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102821\  
 Data File : BP007756.D  
 Acq On : 28 Oct 2021 16:45  
 Operator : CG/JU  
 Sample : M4282-11  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BFY81

Quant Time: Oct 28 17:10:50 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP102521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 27 15:37:43 2021  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.928  | 152  | 161391   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.734 | 136  | 667448   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.569 | 164  | 452257   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.322 | 188  | 975877   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.410 | 240  | 1068058  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 23.851 | 264  | 1137692  | 20.000 | ng/u1 | -0.02    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.370  | 96  | 18239   | 5.134  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.817  | 84  | 60293   | 5.612  | ng/u1 | 0.04  |
| 7) Phenol-d5                  | 7.093  | 99  | 190237  | 15.519 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.258  | 67  | 267450  | 39.297 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.463  | 132 | 337409  | 35.389 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.640  | 113 | 275351  | 28.686 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 9.087  | 128 | 194008  | 42.653 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.810  | 143 | 212159  | 42.707 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.352 | 165 | 375517  | 37.136 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.869 | 131 | 454945  | 34.026 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.975 | 166 | 1303527 | 42.551 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.257 | 160 | 1511709 | 40.261 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.757 | 143 | 79046   | 14.076 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.563 | 176 | 1113330 | 43.109 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.675 | 200 | 197194  | 37.777 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.422 | 188 | 1976024 | 49.272 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.657 | 212 | 2401812 | 48.301 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.698 | 264 | 2578026 | 46.128 | ng/u1 | -0.01 |

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

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

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