

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090823\  
 Data File : BP017020.D  
 Acq On : 08 Sep 2023 19:40  
 Operator : MA/JU  
 Sample : 04175-23  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EZYP0

Quant Time: Sep 09 01:20:41 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP090523.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 06 01:52:05 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.975  | 152  | 394303   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.799 | 136  | 1734297  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10      | 14.628 | 164  | 1042733  | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10      | 17.398 | 188  | 2188921  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12          | 21.492 | 240  | 1651379  | 20.000 | ng/u1 | #-0.02   |
| 88) Perylene-d12          | 24.027 | 264  | 1575045  | 20.000 | ng/u1 | -0.02    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.323  | 96  | 53748   | 5.156  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.764  | 84  | 351781  | 11.810 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.122  | 99  | 335356  | 9.309  | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.317  | 67  | 972383  | 43.154 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.493  | 132 | 860723  | 33.014 | ng/u1 | -0.01 |
| 15) 4-Methylphenol-d8         | 8.681  | 113 | 596285  | 21.727 | ng/u1 | -0.01 |
| 21) Nitrobenzene-d5           | 9.175  | 128 | 559256  | 41.565 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.887  | 143 | 570794  | 39.135 | ng/u1 | -0.01 |
| 28) 2,4-Dichlorophenol-d3     | 10.410 | 165 | 947745  | 35.197 | ng/u1 | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.963 | 131 | 1348045 | 34.700 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 14.046 | 166 | 3475195 | 44.773 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.334 | 160 | 3766935 | 42.053 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.846 | 143 | 109742  | 7.238  | ng/u1 | -0.01 |
| 60) Fluorene-d10              | 15.628 | 176 | 2923441 | 43.991 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.751 | 200 | 419898  | 31.876 | ng/u1 | -0.01 |
| 73) Anthracene-d10            | 17.498 | 188 | 4641538 | 46.173 | ng/u1 | -0.01 |
| 81) Pyrene-d10                | 19.734 | 212 | 5082011 | 50.524 | ng/u1 | -0.01 |
| 92) Benzo(a)pyrene-d12        | 23.863 | 264 | 3916109 | 49.010 | ng/u1 | -0.02 |

Target Compounds Qvalue

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

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