

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070124\  
 Data File : BP020739.D  
 Acq On : 02 Jul 2024 06:35  
 Operator : MA/JU  
 Sample : P2962-11  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 COT83

Quant Time: Jul 02 07:02:50 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062524.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jun 25 18:30:06 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.740  | 152  | 279362   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.504 | 136  | 1245418  | 20.000 | ng/u1 | -0.02    |
| 38) Acenaphthene-d10          | 14.375 | 164  | 840815   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.180 | 188  | 1742900  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.645 | 240  | 1203401  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 25.004 | 264  | 1305746  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.269  | 96   | 20839    | 3.203  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 117871   | 6.217  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.922  | 99   | 465170   | 20.302 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.081  | 67   | 278134   | 19.148 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.281  | 132  | 379644   | 20.273 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.440  | 113  | 400270   | 21.509 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.887  | 128  | 198183   | 21.554 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.599  | 143  | 225124   | 24.466 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.134 | 165  | 410131   | 21.638 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.640 | 131  | 524925   | 19.762 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.781 | 166  | 1406417  | 24.057 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.063 | 160  | 1569764  | 22.503 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.592 | 143  | 188943   | 21.153 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.381 | 176  | 1211138  | 24.619 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.516 | 200  | 163940   | 18.665 | ng/u1 | 0.01     |
| 73) Anthracene-d10            | 17.280 | 188  | 1888777  | 24.146 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.669 | 212  | 1967047  | 27.208 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.786 | 264  | 1583962  | 24.938 | ng/u1 | 0.02     |

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

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

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