

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062022\  
 Data File : BP010768.D  
 Acq On : 20 Jun 2022 19:06  
 Operator : CG/JU  
 Sample : N3358-05  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EW4S1

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :mohammad ahmed 06/27/2022

Quant Time: Jun 21 00:46:37 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 20 16:18:42 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.816  | 152  | 121828   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.610 | 136  | 457088   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 273708   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.210 | 188  | 545883   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.310 | 240  | 574232   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.698 | 264  | 658111   | 20.000 | ng/u1 | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.275  | 96   | 17569    | 6.198  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.699  | 84   | 77123    | 10.462 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.004  | 99   | 66125    | 7.480  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.151  | 67   | 192818   | 35.824 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.351  | 132  | 215053   | 28.583 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.540  | 113  | 127908   | 17.907 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.981  | 128  | 127619m  | 37.604 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.704  | 143  | 132706   | 36.498 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.251 | 165  | 209587   | 31.110 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.763 | 131  | 270442   | 28.499 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.869 | 166  | 768941   | 40.919 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.145 | 160  | 883729   | 38.832 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.733 | 143  | 14549m   | 5.241  | ng/u1 | 0.05     |
| 60) Fluorene-d10              | 15.451 | 176  | 685549   | 41.042 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.592 | 200  | 101921   | 31.294 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.310 | 188  | 1108543m | 44.779 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.551 | 212  | 1337771  | 46.863 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.545 | 264  | 1504988  | 45.925 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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