

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP041223\  
 Data File : BP014673.D  
 Acq On : 12 Apr 2023 14:56  
 Operator : CG/JU  
 Sample : 02251-15  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 COMD3

Quant Time: Apr 13 00:49:30 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP032823.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Apr 11 01:21:43 2023  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

Reviewed By :Christian Giraldo 04/13/2023  
 Supervised By :Jagrut Upadhyay 04/13/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.981  | 152  | 110462   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.799 | 136  | 466659   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.628 | 164  | 254284   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.392 | 188  | 480432   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.486 | 240  | 322999   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.992 | 264  | 433992   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.364  | 96   | 12876    | 4.520  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.805  | 84   | 74386    | 10.190 | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 7.158  | 99   | 31140    | 3.057  | ng/u1 | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.305  | 67   | 190613   | 29.162 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.511  | 132  | 97555    | 13.249 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.705  | 113  | 79097    | 9.626  | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.158  | 128  | 93848    | 24.911 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.881  | 143  | 61226    | 14.193 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.428 | 165  | 100253   | 12.761 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.952 | 131  | 166410   | 16.081 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.040 | 166  | 580909   | 28.175 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.322 | 160  | 630089   | 27.542 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.016 | 143  | 3821m    | 1.227  | ng/u1 | 0.16     |
| 60) Fluorene-d10              | 15.628 | 176  | 472005   | 27.452 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.769 | 200  | 30453    | 8.614  | ng/u1 | 0.01     |
| 73) Anthracene-d10            | 17.492 | 188  | 685105   | 29.466 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.728 | 212  | 641917   | 29.587 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.821 | 264  | 701813   | 30.437 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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