

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP041522\  
 Data File : BP009871.D  
 Acq On : 15 Apr 2022 21:14  
 Operator : CG/JU  
 Sample : PB144125BL  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK125

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 04/18/2022  
 Supervised By :Yogesh Patel 04/18/2022

Quant Time: Apr 18 01:04:18 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP040722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Apr 18 00:59:30 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.958  | 152  | 121773   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.763 | 136  | 562139   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.587 | 164  | 414407   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.339 | 188  | 949537   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.422 | 240  | 979446   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.886 | 264  | 877061   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.381  | 96   | 18790    | 6.887  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.799  | 84   | 266178   | 35.382 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.111  | 99   | 345874   | 32.352 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.281  | 67   | 238671   | 37.499 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.487  | 132  | 285926   | 35.456 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.658  | 113  | 310521   | 34.794 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.116  | 128  | 157454   | 38.842 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.840  | 143  | 174639   | 39.295 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.381 | 165  | 303397   | 32.623 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.893 | 131  | 463439   | 34.987 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.993 | 166  | 1204874  | 37.014 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.281 | 160  | 1298648  | 33.674 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.769 | 143  | 197123   | 33.391 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 1004613  | 36.686 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.692 | 200  | 182845m  | 32.505 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.439 | 188  | 1690689  | 37.188 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.669 | 212  | 2081136  | 38.905 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.727 | 264  | 1829907  | 40.064 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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