

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM090122\  
 Data File : BM036475.D  
 Acq On : 01 Sep 2022 16:23  
 Operator : CG/JU  
 Sample : PB147416BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK416

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 09/02/2022  
 Supervised By :Sohil Jodhani 09/02/2022

Quant Time: Sep 02 02:21:36 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 02 01:33:23 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.722  | 152  | 187397   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.522 | 136  | 853090   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.374 | 164  | 528627   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.121 | 188  | 1099231  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.309 | 240  | 1112552  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.609 | 264  | 1132543  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.128  | 96   | 30385    | 6.179  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.551  | 84   | 395446   | 29.163 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.898  | 99   | 503605   | 31.483 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.057  | 67   | 357462   | 35.502 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.251  | 132  | 438195   | 34.866 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.445  | 113  | 422438   | 35.370 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.892  | 128  | 218167   | 33.730 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.610  | 143  | 222641   | 33.935 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.145 | 165  | 379414   | 32.365 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.675 | 131  | 341167   | 18.773 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.798 | 166  | 1306718  | 39.073 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.068 | 160  | 1552411  | 37.217 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.598 | 143  | 219325m  | 34.611 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.368 | 176  | 1166761  | 38.609 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.504 | 200  | 174876   | 30.944 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.221 | 188  | 1854895  | 37.950 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.521 | 212  | 2196724  | 38.063 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.462 | 264  | 2125807  | 38.283 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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