

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072122\  
 Data File : BP011074.D  
 Acq On : 21 Jul 2022 22:31  
 Operator : CG/JU  
 Sample : PB146324BL  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Jul 21 23:26:26 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071422.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 19 03:25:54 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.669  | 152  | 467728   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.463 | 136  | 1792757  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.334 | 164  | 910903   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.098 | 188  | 1667404  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.222 | 240  | 1582020  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.569 | 264  | 1697579  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.176  | 96   | 84464    | 6.510  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.581  | 84   | 1020907  | 30.443 | ng/u1 | -0.01    |
| 7) Phenol-d5                  | 6.846  | 99   | 1197641  | 31.193 | ng/u1 | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.017  | 67   | 789820   | 31.991 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.205  | 132  | 1011737  | 32.612 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.387  | 113  | 916522   | 29.818 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.834  | 128  | 455251   | 34.846 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.552  | 143  | 450160   | 34.822 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.087 | 165  | 739493   | 31.279 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.610 | 131  | 1136059  | 29.148 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.751 | 166  | 2121643  | 34.079 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.022 | 160  | 2645882  | 36.020 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.546 | 143  | 299730   | 23.788 | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.334 | 176  | 1739067  | 32.863 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.457 | 200  | 216534   | 26.244 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.198 | 188  | 2534240  | 35.751 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.457 | 212  | 2795633  | 33.647 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.416 | 264  | 2973984  | 36.476 | ng/u1 | -0.01    |

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

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

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