

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM081422\  
 Data File : BM036271.D  
 Acq On : 14 Aug 2022 15:31  
 Operator : CG/JU  
 Sample : N4166-05  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :

Quant Time: Aug 14 23:37:59 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM072522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Aug 14 23:34:05 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.810  | 152  | 247196   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.610 | 136  | 1065551  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.451 | 164  | 627013   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.192 | 188  | 1214310  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.374 | 240  | 1082804  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.709 | 264  | 1100067  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.210  | 96   | 25434    | 3.872  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.652  | 84   | 73160    | 4.065  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 6.998  | 99   | 71282    | 3.439  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.145  | 67   | 303529   | 21.964 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.345  | 132  | 264500   | 15.372 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.539  | 113  | 150107   | 8.938  | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.981  | 128  | 196672   | 23.293 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.698  | 143  | 187793   | 22.828 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.239 | 165  | 292130   | 19.635 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.763 | 131  | 499753   | 19.842 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.869 | 166  | 1178076  | 27.583 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.145 | 160  | 1373233  | 25.199 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.704 | 143  | 5183     | 0.591  | ng/u1 | 0.02     |
| 60) Fluorene-d10              | 15.445 | 176  | 1062398  | 27.527 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.574 | 200  | 126245   | 19.341 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.292 | 188  | 1683026  | 30.656 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.586 | 212  | 1823965  | 30.394 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.556 | 264  | 1695704  | 30.362 | ng/u1 | 0.00     |

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

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

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