

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM072722\  
 Data File : BM035982.D  
 Acq On : 28 Jul 2022 13:10  
 Operator : CG/JU  
 Sample : N3874-07  
 Misc :  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 GBJ31

Quant Time: Jul 29 00:49:29 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM072522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jul 27 23:55:50 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.875  | 152  | 300401   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8        | 10.681 | 136  | 1239896  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10      | 14.516 | 164  | 677368   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.257 | 188  | 1224719  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.433 | 240  | 1117751  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.809 | 264  | 1176239  | 20.000 | ng/u1 | 0.00     |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.269  | 96  | 37885   | 4.746  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.693  | 84  | 184841  | 8.452  | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.040  | 99  | 125099  | 4.966  | ng/u1 | -0.01 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.204  | 67  | 302501  | 18.013 | ng/u1 | -0.01 |
| 11) 2-Chlorophenol-d4         | 7.404  | 132 | 320947  | 15.349 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.587  | 113 | 209142  | 10.248 | ng/u1 | -0.01 |
| 21) Nitrobenzene-d5           | 9.040  | 128 | 195590  | 19.907 | ng/u1 | -0.01 |
| 24) 2-Nitrophenol-d4          | 9.763  | 143 | 183262  | 19.145 | ng/u1 | -0.01 |
| 28) 2,4-Dichlorophenol-d3     | 10.298 | 165 | 300744  | 17.371 | ng/u1 | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.822 | 131 | 577873  | 19.717 | ng/u1 | -0.01 |
| 46) Dimethylphthalate-d6      | 13.928 | 166 | 985595  | 21.361 | ng/u1 | -0.01 |
| 49) Acenaphthylene-d8         | 14.210 | 160 | 1204655 | 20.463 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.716 | 143 | 35678   | 3.765  | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.504 | 176 | 835362  | 20.035 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.622 | 200 | 96993   | 14.733 | ng/u1 | -0.01 |
| 73) Anthracene-d10            | 17.351 | 188 | 1242524 | 22.440 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.645 | 212 | 1388953 | 22.421 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.656 | 264 | 1316432 | 22.045 | ng/u1 | 0.00  |

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

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

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