

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM063021\  
 Data File : BM030722.D  
 Acq On : 01 Jul 2021 01:35  
 Operator : CG/JU  
 Sample : M2808-03  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BGG09

Manual Integrations  
 APPROVED

mohammad  
 7/1/2021 2:55:29 PM

Quant Time: Jul 01 02:46:14 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM062221.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jul 01 00:59:58 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min)     |
|-------------------------------|--------|------|----------|--------|-------|--------------|
| -----                         |        |      |          |        |       |              |
| Internal Standards            |        |      |          |        |       |              |
| 1) 1,4-Dichlorobenzene-d4     | 7.775  | 152  | 81422    | 20.000 | ng/u1 | 0.00         |
| 20) Naphthalene-d8            | 10.563 | 136  | 333286   | 20.000 | ng/u1 | 0.00         |
| 38) Acenaphthene-d10          | 14.404 | 164  | 217927   | 20.000 | ng/u1 | 0.00         |
| 64) Phenanthrene-d10          | 17.145 | 188  | 444746   | 20.000 | ng/u1 | 0.00         |
| 79) Chrysene-d12              | 21.309 | 240  | 381377   | 20.000 | ng/u1 | 0.00         |
| 88) Perylene-d12              | 23.562 | 264  | 356792   | 20.000 | ng/u1 | 0.00         |
| System Monitoring Compounds   |        |      |          |        |       |              |
| 3) 1,4-Dioxane-d8             | 3.269  | 96   | 4447m    | 2.214  | ng/uL | 0.00         |
| 4) Pyridine-d5                | 3.716  | 84   | 20130m   | 3.731  | ng/u1 | 0.04         |
| 7) Phenol-d5                  | 6.946  | 99   | 67425    | 9.602  | ng/u1 | 0.00         |
| 9) Bis-(2-Chloroethyl)eth...  | 7.110  | 67   | 48142    | 10.907 | ng/u1 | 0.00         |
| 11) 2-Chlorophenol-d4         | 7.310  | 132  | 58679    | 10.826 | ng/u1 | 0.00         |
| 15) 4-Methylphenol-d8         | 8.481  | 113  | 53069    | 9.711  | ng/u1 | 0.00         |
| 21) Nitrobenzene-d5           | 8.934  | 128  | 25651    | 10.314 | ng/u1 | 0.00         |
| 24) 2-Nitrophenol-d4          | 9.651  | 143  | 26665    | 9.648  | ng/u1 | 0.00         |
| 28) 2,4-Dichlorophenol-d3     | 10.187 | 165  | 50279    | 9.496  | ng/u1 | 0.00         |
| 31) 4-Chloroaniline-d4        | 10.704 | 131  | 66084    | 8.950  | ng/u1 | 0.00         |
| 46) Dimethylphthalate-d6      | 13.816 | 166  | 168280   | 11.183 | ng/u1 | 0.00         |
| 49) Acenaphthylene-d8         | 14.098 | 160  | 201069   | 10.294 | ng/u1 | 0.00         |
| 54) 4-Nitrophenol-d4          | 14.627 | 143  | 10511m   | 3.687  | ng/u1 | 0.02         |
| 60) Fluorene-d10              | 15.398 | 176  | 151466   | 11.349 | ng/u1 | 0.00         |
| 65) 4,6-Dinitro-2-methylph... | 15.527 | 200  | 10590    | 3.639  | ng/u1 | 0.00         |
| 73) Anthracene-d10            | 17.245 | 188  | 223682   | 10.777 | ng/u1 | 0.00         |
| 81) Pyrene-d10                | 19.527 | 212  | 231986   | 11.609 | ng/u1 | 0.00         |
| 92) Benzo(a)pyrene-d12        | 23.421 | 264  | 170891   | 9.213  | ng/u1 | 0.00         |
| Target Compounds              |        |      |          |        |       |              |
| 47) Dimethylphthalate         | 13.863 | 163  | 27615    | 1.873  | ng/u1 | Qvalue<br>99 |
| -----                         |        |      |          |        |       |              |

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

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