

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051021\  
 Data File : BM029913.D  
 Acq On : 10 May 2021 20:11  
 Operator : CG/JU  
 Sample : M2177-02  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BG579

Manual Integrations  
 APPROVED

mohammad  
 5/11/2021 2:33:02 PM

Quant Time: May 11 01:39:10 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050721.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 11 01:19:00 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.55  | 152  | 135076   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.33 | 136  | 621380   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.20 | 164  | 412698   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 16.94 | 188  | 848951   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.13 | 240  | 837528   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.29 | 264  | 751533   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.07  | 96  | 4003   | 1.18  | ng/uL | 0.01 |
| 4) Pyridine-d5                 | 3.55  | 84  | 28152m | 3.32  | ng/ul | 0.06 |
| 7) Phenol-d5                   | 6.75  | 99  | 92128  | 7.33  | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 6.90  | 67  | 41526  | 5.33  | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.09  | 132 | 62566  | 6.45  | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.28  | 113 | 43611  | 4.20  | ng/ul | 0.01 |
| 21) Nitrobenzene-d5            | 8.72  | 128 | 20810  | 4.86  | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.43  | 143 | 22003  | 4.55  | ng/ul | 0.01 |
| 28) 2,4-Dichlorophenol-d3      | 9.97  | 165 | 58524  | 6.04  | ng/ul | 0.02 |
| 31) 4-Chloroaniline-d4         | 10.49 | 131 | 53164  | 3.60  | ng/ul | 0.02 |
| 46) Dimethylphthalate-d6       | 13.62 | 166 | 201533 | 6.30  | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 13.89 | 160 | 410857 | 10.13 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.49 | 143 | 8967m  | 1.72  | ng/ul | 0.06 |
| 60) Fluorene-d10               | 15.20 | 176 | 186941 | 6.88  | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.34 | 200 | 14032  | 2.54  | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.04 | 188 | 452390 | 10.58 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.34 | 212 | 951705 | 22.15 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.15 | 264 | 858241 | 20.15 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |              | Ovalue |
|--------------------------------|-------|-----|--------|--------------|--------|
| 72) Phenanthrene               | 16.99 | 178 | 50541  | 1.010 ng/ul  | 99     |
| 80) Fluoranthene               | 19.01 | 202 | 139941 | 2.676 ng/ul  | 98     |
| 82) Pyrene                     | 19.37 | 202 | 292523 | 5.316 ng/ul  | 100    |
| 85) Benzo(a)anthracene         | 21.12 | 228 | 86398  | 1.581 ng/ul  | 94     |
| 86) Bis(2-ethylhexyl)phthalate | 21.06 | 149 | 68724  | 2.038 ng/ul  | 99     |
| 87) Chrysene                   | 21.17 | 228 | 99535  | 1.829 ng/ul  | 98     |
| 90) Benzo(b)fluoranthene       | 22.65 | 252 | 95005m | 1.949 ng/ul  |        |
| 93) Benzo(a)pyrene             | 23.20 | 252 | 91356  | 2.063 ng/ul# | 95     |
| 96) Benzo(a,h,i)perylene       | 26.10 | 276 | 50270  | 1.158 ng/ul  | 96     |

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

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