

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021020\  
 Data File : BM024808.D  
 Acq On : 10 Feb 2020 15:04  
 Operator : CG/JU  
 Sample : L1402-03  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 CB6L1

Manual Integrations  
 APPROVED

mohammad  
 2/12/2020 9:50:15 AM

Quant Time: Feb 10 16:47:12 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM020620MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Feb 10 16:33:34 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.40  | 152  | 274469   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.16 | 136  | 1078053  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.06 | 164  | 775146   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.81 | 188  | 1680911  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.02 | 240  | 1667758  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.13 | 264  | 1754155  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.02  | 96  | 20072   | 3.52  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.60  | 99  | 336207  | 18.44 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.75  | 67  | 203927  | 19.80 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 6.95  | 132 | 309245  | 18.93 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.12  | 113 | 251531  | 16.35 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.54  | 128 | 149114  | 19.13 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.26  | 143 | 173195  | 18.80 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.79  | 165 | 319980  | 17.22 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.30 | 131 | 362974  | 20.88 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.48 | 166 | 1020071 | 17.43 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.74 | 160 | 1205120 | 17.64 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.28 | 143 | 114270  | 12.02 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.06 | 176 | 883057  | 17.04 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.19 | 200 | 135533  | 12.37 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.90 | 188 | 1273157 | 16.59 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.22 | 212 | 1500548 | 17.99 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 22.99 | 264 | 1564945 | 17.76 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |              | Ovalue |
|----------------------------|-------|-----|---------|--------------|--------|
| 70) Phenanthrene           | 16.85 | 178 | 339648  | 3.706 ng/ul  | 98     |
| 77) Fluoranthene           | 18.88 | 202 | 613246  | 5.451 ng/ul  | 99     |
| 80) Pyrene                 | 19.24 | 202 | 521043  | 4.720 ng/ul  | 99     |
| 83) Benzo(a)anthracene     | 21.01 | 228 | 293901  | 2.616 ng/ul  | 96     |
| 85) Chrysene               | 21.06 | 228 | 289666  | 2.619 ng/ul  | 100    |
| 88) Benzo(b)fluoranthene   | 22.51 | 252 | 412843  | 3.701 ng/ul# | 97     |
| 89) Benzo(k)fluoranthene   | 22.54 | 252 | 127242m | 1.185 ng/ul  |        |
| 91) Benzo(a)pyrene         | 23.03 | 252 | 249048  | 2.490 ng/ul  | 99     |
| 92) Indeno(1,2,3-cd)pyrene | 25.17 | 276 | 184009  | 1.478 ng/ul  | 98     |
| 94) Benzo(g,h,i)perylene   | 25.79 | 276 | 155941  | 1.506 ng/ul  | 98     |

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

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