

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021920\  
 Data File : BM024948.D  
 Acq On : 19 Feb 2020 17:40  
 Operator : CG/JU  
 Sample : L1486-15DL 10X  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBCW6DL

Manual Integrations  
 APPROVED

mohammad  
 2/20/2020 3:12:37 PM

Quant Time: Feb 19 23:53:17 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM020620MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Feb 19 03:36:08 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.39  | 152  | 409981   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.14 | 136  | 1516964  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.03 | 164  | 929859   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.79 | 188  | 2051023  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.00 | 240  | 2198916  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.10 | 264  | 2281735  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |      |       |       |
|--------------------------------|-------|-----|--------|------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.05  | 96  | 1219m  | 0.14 | ng/uL | 0.01  |
| 5) Phenol-d5                   | 6.59  | 99  | 14930  | 0.55 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.75  | 67  | 44766  | 2.91 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 6.93  | 132 | 48462  | 1.99 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.10  | 113 | 26860  | 1.17 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 8.53  | 128 | 30119  | 2.75 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 9.24  | 143 | 28965  | 2.23 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 9.77  | 165 | 55403  | 2.12 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 0.00  | 131 | 0d     | 0.00 | ng/ul |       |
| 44) Dimethylphthalate-d6       | 13.46 | 166 | 191425 | 2.73 | ng/ul | 0.00  |
| 47) Acenaphthylene-d8          | 13.72 | 160 | 229940 | 2.81 | ng/ul | 0.00  |
| 52) 4-Nitrophenol-d4           | 0.00  | 143 | 0d     | 0.00 | ng/ul |       |
| 58) Fluorene-d10               | 15.04 | 176 | 175462 | 2.82 | ng/ul | 0.00  |
| 63) 4,6-Dinitro-2-methylphenol | 15.17 | 200 | 16773  | 1.25 | ng/ul | 0.00  |
| 71) Anthracene-d10             | 16.89 | 188 | 270517 | 2.89 | ng/ul | -0.01 |
| 79) Pyrene-d10                 | 19.20 | 212 | 306107 | 2.78 | ng/ul | 0.00  |
| 90) Benzo(a)pyrene-d12         | 22.97 | 264 | 321488 | 2.81 | ng/ul | 0.00  |

## Target Compounds

|                         |       |     |           | Ovalue  |        |
|-------------------------|-------|-----|-----------|---------|--------|
| 24) 2,4-Dimethylphenol  | 9.35  | 107 | 42260m    | 1.621   | ng/ul  |
| 28) Naphthalene         | 10.19 | 128 | 19600009m | 252.878 | ng/ul  |
| 34) 2-Methylnaphthalene | 11.82 | 142 | 2894851   | 50.703  | ng/ul  |
| 41) 1,1'-Biphenyl       | 12.86 | 154 | 786271    | 11.454  | ng/ul# |
| 50) Acenaphthene        | 14.10 | 153 | 3519636   | 61.382  | ng/ul  |
| 54) Dibenzofuran        | 14.44 | 168 | 2959529   | 34.625  | ng/ul  |
| 59) Fluorene            | 15.09 | 166 | 1380356   | 19.595  | ng/ul  |
| 70) Phenanthrene        | 16.83 | 178 | 2918059   | 26.093  | ng/ul  |
| 75) Carbazole           | 17.20 | 167 | 785246    | 8.311   | ng/ul  |
| 77) Fluoranthene        | 18.86 | 202 | 402152    | 2.929   | ng/ul# |
| 80) Pyrene              | 19.22 | 202 | 255220    | 1.753   | ng/ul# |

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

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