

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\  
 Data File : BM026989.D  
 Acq On : 03 Aug 2020 21:40  
 Operator : JU/CG  
 Sample : L3424-03DL 10X  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0S78DL

Manual Integrations  
 APPROVED

mohammad  
 8/5/2020 10:26:35 AM

Quant Time: Aug 04 04:53:02 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM072720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 03 10:21:18 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.66  | 152  | 104998   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.43 | 136  | 412423   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.29 | 164  | 249153   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.03 | 188  | 522075   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.22 | 240  | 523390   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.43 | 264  | 519012   | 20.00 | ng/ul | 0.00     |

|                                |       |     |       |      |       |      |
|--------------------------------|-------|-----|-------|------|-------|------|
| System Monitoring Compounds    |       |     |       |      |       |      |
| 3) 1,4-Dioxane-d8              | 3.22  | 96  | 447   | 0.19 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.84  | 99  | 13778 | 1.50 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.00  | 67  | 10134 | 1.63 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.20  | 132 | 10766 | 1.56 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.36  | 113 | 10832 | 1.53 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.80  | 128 | 5057  | 1.58 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.52  | 143 | 4962  | 1.69 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.06 | 165 | 10962 | 1.56 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.57 | 131 | 6201  | 0.74 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.70 | 166 | 35088 | 1.80 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.98 | 160 | 40881 | 1.64 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.49 | 143 | 2323  | 0.70 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.29 | 176 | 28462 | 1.67 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.40 | 200 | 2112  | 0.91 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.13 | 188 | 45515 | 1.74 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.43 | 212 | 49283 | 1.77 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.29 | 264 | 48201 | 1.65 | ng/ul | 0.00 |

|                            |       |     |         |        |        |     |
|----------------------------|-------|-----|---------|--------|--------|-----|
| Target Compounds           |       |     |         | Ovalue |        |     |
| 28) Naphthalene            | 10.48 | 128 | 51640   | 2.241  | ng/ul  | 97  |
| 48) Acenaphthylene         | 14.01 | 152 | 138972  | 5.502  | ng/ul  | 98  |
| 70) Phenanthrene           | 17.07 | 178 | 84838   | 2.677  | ng/ul  | 98  |
| 72) Anthracene             | 17.16 | 178 | 453672  | 13.928 | ng/ul  | 99  |
| 75) Carbazole              | 17.43 | 167 | 33136   | 1.151  | ng/ul# | 97  |
| 77) Fluoranthene           | 19.09 | 202 | 566635  | 15.143 | ng/ul  | 100 |
| 80) Pyrene                 | 19.46 | 202 | 680324  | 17.866 | ng/ul  | 97  |
| 83) Benzo(a)anthracene     | 21.20 | 228 | 475024  | 13.168 | ng/ul  | 94  |
| 85) Chrysene               | 21.26 | 228 | 850040  | 24.163 | ng/ul  | 99  |
| 88) Benzo(b)fluoranthene   | 22.78 | 252 | 1824346 | 49.137 | ng/ul  | 99  |
| 89) Benzo(k)fluoranthene   | 22.82 | 252 | 572728m | 16.076 | ng/ul  |     |
| 91) Benzo(a)pyrene         | 23.33 | 252 | 526926  | 15.728 | ng/ul  | 99  |
| 92) Indeno(1,2,3-cd)pyrene | 25.64 | 276 | 739615  | 17.795 | ng/ul  | 95  |
| 93) Dibenzo(a,h)anthracene | 25.64 | 278 | 189367m | 5.387  | ng/ul  |     |
| 94) Benzo(g,h,i)perylene   | 26.31 | 276 | 351020  | 10.214 | ng/ul  | 97  |

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

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