

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081320\  
 Data File : BM027098.D  
 Acq On : 14 Aug 2020 22:08  
 Operator : JU/CG  
 Sample : L3631-01  
 Misc :  
 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BFML7

Manual Integrations  
 APPROVED

mohammad  
 8/19/2020 8:18:50 AM

Quant Time: Aug 17 07:40:42 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM080720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 13 11:47:07 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.64  | 152  | 94702    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.41 | 136  | 332502   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.27 | 164  | 212654   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.02 | 188  | 470367m  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.21 | 240  | 603657   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.40 | 264  | 669840   | 20.00 | ng/ul | 0.00     |

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| System Monitoring Compounds    |       |     |        |       |       |      |
| 3) 1,4-Dioxane-d8              | 3.20  | 96  | 5206   | 2.50  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.82  | 99  | 99491  | 13.51 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.97  | 67  | 75535  | 14.97 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.17  | 132 | 82668  | 14.23 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.34  | 113 | 70875  | 12.35 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.78  | 128 | 40456  | 15.35 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.50  | 143 | 41913  | 15.35 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.03 | 165 | 87287  | 15.20 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.54 | 131 | 91117  | 13.66 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.69 | 166 | 295229 | 17.41 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.96 | 160 | 309367 | 15.31 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.47 | 143 | 30477  | 10.65 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.27 | 176 | 236935 | 16.03 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.39 | 200 | 27102  | 10.55 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.11 | 188 | 357130 | 15.41 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.41 | 212 | 433957 | 14.41 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.26 | 264 | 562050 | 15.03 | ng/ul | 0.00 |

|                            |       |     |         |        |       |     |
|----------------------------|-------|-----|---------|--------|-------|-----|
| Target Compounds           |       |     |         | Ovalue |       |     |
| 45) Dimethylphthalate      | 13.73 | 163 | 83474   | 4.697  | ng/ul | 100 |
| 70) Phenanthrene           | 17.06 | 178 | 129825m | 4.604  | ng/ul |     |
| 77) Fluoranthene           | 19.07 | 202 | 392515m | 11.479 | ng/ul |     |
| 80) Pyrene                 | 19.44 | 202 | 360376  | 8.795  | ng/ul | 96  |
| 83) Benzo(a)anthracene     | 21.19 | 228 | 215167  | 5.295  | ng/ul | 94  |
| 85) Chrysene               | 21.24 | 228 | 198115  | 4.967  | ng/ul | 97  |
| 88) Benzo(b)fluoranthene   | 22.75 | 252 | 305774  | 6.574  | ng/ul | 98  |
| 89) Benzo(k)fluoranthene   | 22.79 | 252 | 90155m  | 1.956  | ng/ul |     |
| 91) Benzo(a)pyrene         | 23.31 | 252 | 197941  | 4.676  | ng/ul | 96  |
| 92) Indeno(1,2,3-cd)pyrene | 25.59 | 276 | 143615  | 2.633  | ng/ul | 95  |
| 94) Benzo(g,h,i)perylene   | 26.26 | 276 | 106548  | 2.365  | ng/ul | 96  |

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

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