

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062519\  
 Data File : BM021197.D  
 Acq On : 26 Jun 2019 18:17  
 Operator : HP/JU  
 Sample : K3501-01  
 Misc :  
 ALS Vial : 46 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C5AB9

Manual Integrations  
 APPROVED

mohammad  
 6/27/2019 8:17:55 AM

Quant Time: Jun 27 00:51:40 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 27 00:42:34 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.76  | 152  | 231635   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.55 | 136  | 1057139  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.40 | 164  | 695266   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.16 | 188  | 1569999  | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.34 | 240  | 1075091  | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 23.62 | 264  | 1168440  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.29  | 96  | 15971   | 3.27  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.93  | 99  | 516632  | 25.62 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.10  | 67  | 295914  | 24.63 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.30  | 132 | 410018  | 25.03 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.47  | 113 | 489051  | 29.06 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.93  | 128 | 245805  | 28.64 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.65  | 143 | 293716  | 30.82 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.17 | 165 | 588879  | 30.39 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.70 | 131 | 661804  | 32.75 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.82 | 166 | 2240442 | 35.66 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.10 | 160 | 2478020 | 32.22 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.62 | 143 | 266500  | 26.08 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.40 | 176 | 1904331 | 34.85 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.53 | 200 | 286097  | 29.64 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.25 | 188 | 2845432 | 34.88 | ng/ul | 0.00 |
| 78) Pyrene-d10                 | 19.55 | 212 | 2707501 | 41.97 | ng/ul | 0.00 |
| 89) Benzo(a)pyrene-d12         | 23.48 | 264 | 2394708 | 34.74 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |        | Ovalue |    |
|--------------------------------|-------|-----|--------|--------|--------|----|
| 44) Dimethylphthalate          | 13.86 | 163 | 865715 | 15.138 | ng/ul  | 98 |
| 69) Phenanthrene               | 17.20 | 178 | 273899 | 3.154  | ng/ul  | 99 |
| 76) Fluoranthene               | 19.22 | 202 | 425985 | 4.540  | ng/ul  | 99 |
| 79) Pyrene                     | 19.58 | 202 | 347586 | 4.573  | ng/ul  | 99 |
| 82) Benzo(a)anthracene         | 21.33 | 228 | 160510 | 2.190  | ng/ul  | 99 |
| 83) Bis(2-ethylhexyl)phthalate | 21.26 | 149 | 41845  | 1.045  | ng/ul  | 97 |
| 84) Chrysene                   | 21.38 | 228 | 173232 | 2.524  | ng/ul  | 98 |
| 87) Benzo(b)fluoranthene       | 22.94 | 252 | 223955 | 3.026  | ng/ul# | 98 |
| 88) Benzo(k)fluoranthene       | 22.98 | 252 | 84197m | 1.182  | ng/ul  |    |
| 90) Benzo(a)pyrene             | 23.52 | 252 | 144789 | 2.078  | ng/ul  | 97 |
| 91) Indeno(1,2,3-cd)pyrene     | 25.91 | 276 | 101984 | 1.243  | ng/ul  | 98 |
| 93) Benzo(g,h,i)perylene       | 26.61 | 276 | 86054  | 1.274  | ng/ul  | 99 |

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

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