

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM123119\  
 Data File : BM024400.D  
 Acq On : 31 Dec 2019 17:40  
 Operator : JU  
 Sample : K6308-01  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0C56

Manual Integrations  
 APPROVED

mohammad  
 1/2/2020 8:41:07 AM

Quant Time: Jan 01 05:37:04 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121719MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 01 04:44:04 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.40  | 152  | 183673   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.17 | 136  | 759911   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.07 | 164  | 441725   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.83 | 188  | 909876   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.05 | 240  | 889322   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.17 | 264  | 1098839  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.00  | 96  | 6533m   | 1.57  | ng/uL | 0.02 |
| 5) Phenol-d5                   | 6.62  | 99  | 77290   | 4.44  | ng/ul | 0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.76  | 67  | 201676  | 19.34 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 6.95  | 132 | 192898  | 15.55 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.14  | 113 | 86384   | 6.08  | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.56  | 128 | 101338  | 18.56 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.28  | 143 | 105128  | 17.51 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.83  | 165 | 190204  | 16.30 | ng/ul | 0.02 |
| 29) 4-Chloroaniline-d4         | 0.00  | 131 | 0d      | 0.00  | ng/ul |      |
| 44) Dimethylphthalate-d6       | 13.50 | 166 | 672251  | 20.58 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.76 | 160 | 764391  | 19.53 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.39 | 143 | 12419m  | 2.12  | ng/ul | 0.06 |
| 58) Fluorene-d10               | 15.07 | 176 | 591534  | 20.66 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.24 | 200 | 61927   | 11.16 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.93 | 188 | 867542  | 20.96 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.24 | 212 | 923684  | 21.80 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.04 | 264 | 1156469 | 21.19 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |        |       | Ovalue |    |
|----------------------------|-------|-----|--------|-------|--------|----|
| 6) Phenol                  | 6.65  | 94  | 54317  | 3.025 | ng/ul# | 90 |
| 45) Dimethylphthalate      | 13.55 | 163 | 65164  | 1.920 | ng/ul  | 98 |
| 77) Fluoranthene           | 18.91 | 202 | 150178 | 2.705 | ng/ul  | 96 |
| 80) Pyrene                 | 19.27 | 202 | 140440 | 2.419 | ng/ul  | 98 |
| 83) Benzo(a)anthracene     | 21.04 | 228 | 73329  | 1.294 | ng/ul  | 96 |
| 85) Chrysene               | 21.09 | 228 | 89518  | 1.612 | ng/ul  | 98 |
| 88) Benzo(b)fluoranthene   | 22.55 | 252 | 156307 | 2.377 | ng/ul# | 94 |
| 91) Benzo(a)pyrene         | 23.08 | 252 | 95872  | 1.612 | ng/ul# | 95 |
| 92) Indeno(1,2,3-cd)pyrene | 25.26 | 276 | 76330  | 1.069 | ng/ul  | 96 |
| 94) Benzo(g,h,i)perylene   | 25.91 | 276 | 86332  | 1.480 | ng/ul  | 96 |

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

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