

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM123119\  
 Data File : BM024406.D  
 Acq On : 31 Dec 2019 21:16  
 Operator : JU  
 Sample : K6314-06  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BFF19

Manual Integrations  
 APPROVED

mohammad  
 1/2/2020 8:41:17 AM

Quant Time: Jan 01 05:55:40 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  | 217496   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.17 | 136  | 909009   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.07 | 164  | 519670   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.83 | 188  | 1058344  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.05 | 240  | 1027932  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.17 | 264  | 1272700  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 2.98  | 96  | 31862   | 6.47  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.61  | 99  | 154785  | 7.52  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.76  | 67  | 370075  | 29.96 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 6.95  | 132 | 376513  | 25.64 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.13  | 113 | 254990  | 15.15 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.56  | 128 | 200147  | 30.64 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.27  | 143 | 220293  | 30.67 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.82  | 165 | 375508  | 26.90 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.37 | 131 | 21142m  | 1.25  | ng/ul | 0.04 |
| 44) Dimethylphthalate-d6       | 13.50 | 166 | 1249711 | 32.52 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.76 | 160 | 1475732 | 32.05 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.37 | 143 | 30770m  | 4.46  | ng/ul | 0.04 |
| 58) Fluorene-d10               | 15.07 | 176 | 1085363 | 32.22 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.23 | 200 | 147026  | 22.77 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.93 | 188 | 1645909 | 34.18 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.24 | 212 | 1830301 | 37.38 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.03 | 264 | 2214368 | 35.03 | ng/ul | 0.00 |

| Target Compounds      |       |     |       |       | Ovalue |    |
|-----------------------|-------|-----|-------|-------|--------|----|
| 45) Dimethylphthalate | 13.55 | 163 | 55248 | 1.383 | ng/ul  | 98 |

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

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