

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102419\  
 Data File : BM023389.D  
 Acq On : 24 Oct 2019 21:04  
 Operator : JU  
 Sample : K5346-08DL 5X  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 ETOM1DL

Manual Integrations  
 APPROVED

mohammad  
 10/26/2019 10:04:17 AM

Quant Time: Oct 24 23:22:58 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM102319MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 23 18:47:16 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.65  | 152  | 524671   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.43 | 136  | 1954945  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.29 | 164  | 1139058  | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.04 | 188  | 2215646  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.24 | 240  | 2186604  | 20.00 | ng/ul | 0.01     |
| 86) Perylene-d12          | 23.44 | 264  | 2761016  | 20.00 | ng/ul | 0.02     |

## System Monitoring Compounds

|                                |       |     |        |      |       |      |
|--------------------------------|-------|-----|--------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.18  | 96  | 9422   | 0.85 | ng/uL | 0.02 |
| 5) Phenol-d5                   | 6.83  | 99  | 161773 | 3.67 | ng/ul | 0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.99  | 67  | 100627 | 3.90 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.18  | 132 | 144904 | 4.21 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.36  | 113 | 118317 | 3.31 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.79  | 128 | 66996  | 4.55 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.52  | 143 | 65959  | 4.04 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.06 | 165 | 133443 | 4.16 | ng/ul | 0.01 |
| 29) 4-Chloroaniline-d4         | 10.56 | 131 | 128338 | 3.48 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.71 | 166 | 439409 | 5.08 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.98 | 160 | 487729 | 4.92 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.51 | 143 | 29154  | 1.95 | ng/ul | 0.02 |
| 58) Fluorene-d10               | 15.29 | 176 | 379839 | 5.19 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.43 | 200 | 5164   | 0.37 | ng/ul | 0.02 |
| 71) Anthracene-d10             | 17.14 | 188 | 620157 | 5.92 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.44 | 212 | 646004 | 5.37 | ng/ul | 0.01 |
| 90) Benzo(a)pyrene-d12         | 23.30 | 264 | 810566 | 5.76 | ng/ul | 0.02 |

## Target Compounds

|                            |       |     |          |        | Ovalue    |
|----------------------------|-------|-----|----------|--------|-----------|
| 45) Dimethylphthalate      | 13.76 | 163 | 125728   | 1.448  | ng/ul 100 |
| 48) Acenaphthylene         | 14.01 | 152 | 368917   | 3.592  | ng/ul 98  |
| 70) Phenanthrene           | 17.08 | 178 | 247454   | 1.998  | ng/ul 99  |
| 72) Anthracene             | 17.17 | 178 | 823850   | 6.567  | ng/ul 99  |
| 77) Fluoranthene           | 19.10 | 202 | 854422   | 6.616  | ng/ul 99  |
| 80) Pyrene                 | 19.47 | 202 | 1296000  | 8.393  | ng/ul 98  |
| 83) Benzo(a)anthracene     | 21.22 | 228 | 1553245  | 10.621 | ng/ul 95  |
| 85) Chrysene               | 21.27 | 228 | 2817235  | 20.879 | ng/ul 99  |
| 88) Benzo(b)fluoranthene   | 22.79 | 252 | 5521878  | 31.139 | ng/ul 98  |
| 89) Benzo(k)fluoranthene   | 22.83 | 252 | 1338257m | 8.037  | ng/ul     |
| 91) Benzo(a)pyrene         | 23.35 | 252 | 3160310  | 19.686 | ng/ul 98  |
| 92) Indeno(1,2,3-cd)pyrene | 25.66 | 276 | 2363390  | 13.596 | ng/ul 96  |
| 93) Dibenzo(a,h)anthracene | 25.66 | 278 | 634930   | 4.332  | ng/ul# 85 |
| 94) Benzo(g,h,i)perylene   | 26.34 | 276 | 1744408  | 12.334 | ng/ul 98  |

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

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