

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN053119\  
 Data File : BN005990.D  
 Acq On : 31 May 2019 21:11  
 Operator : JU/SJ  
 Sample : K2924-22  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 A42F4

Manual Integrations  
 APPROVED

mohammad  
 6/5/2019 6:11:18 AM

Quant Time: Jun 01 02:36:25 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN052819MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jun 01 01:49:29 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.75  | 152  | 542114   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.54 | 136  | 2269722  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.41 | 164  | 1177555  | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.16 | 188  | 2065084  | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.38 | 240  | 1277197  | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 23.68 | 264  | 1510692  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.27  | 96  | 52748   | 4.28  | ng/uL | 0.01 |
| 5) Phenol-d5                   | 6.92  | 99  | 1105515 | 23.01 | ng/ul | 0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.08  | 67  | 661732  | 24.14 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.28  | 132 | 908614  | 24.82 | ng/ul | 0.01 |
| 13) 4-Methylphenol-d8          | 8.46  | 113 | 771019  | 19.88 | ng/ul | 0.01 |
| 19) Nitrobenzene-d5            | 8.91  | 128 | 461054  | 31.65 | ng/ul | 0.01 |
| 22) 2-Nitrophenol-d4           | 9.62  | 143 | 487148  | 35.91 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.17 | 165 | 864121  | 26.97 | ng/ul | 0.01 |
| 29) 4-Chloroaniline-d4         | 10.68 | 131 | 763158  | 18.17 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.81 | 166 | 2454250 | 28.27 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.09 | 160 | 2982573 | 27.47 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.62 | 143 | 303435  | 19.21 | ng/ul | 0.01 |
| 57) Fluorene-d10               | 15.41 | 176 | 1935474 | 26.59 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.54 | 200 | 44713   | 5.31  | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.26 | 188 | 2333399 | 26.55 | ng/ul | 0.00 |
| 78) Pyrene-d10                 | 19.57 | 212 | 2017854 | 31.28 | ng/ul | 0.00 |
| 89) Benzo(a)pyrene-d12         | 23.53 | 264 | 1798538 | 26.68 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |       | Ovalue |     |
|----------------------------|-------|-----|---------|-------|--------|-----|
| 44) Dimethylphthalate      | 13.86 | 163 | 766570  | 8.835 | ng/ul  | 100 |
| 69) Phenanthrene           | 17.21 | 178 | 444111  | 4.341 | ng/ul  | 98  |
| 71) Anthracene             | 17.29 | 178 | 105201  | 1.013 | ng/ul  | 97  |
| 76) Fluoranthene           | 19.23 | 202 | 745722  | 6.827 | ng/ul  | 96  |
| 79) Pyrene                 | 19.60 | 202 | 610750  | 7.549 | ng/ul  | 99  |
| 82) Benzo(a)anthracene     | 21.36 | 228 | 308555  | 3.843 | ng/ul  | 92  |
| 84) Chrysene               | 21.41 | 228 | 310926  | 4.152 | ng/ul# | 93  |
| 87) Benzo(b)fluoranthene   | 22.98 | 252 | 472555  | 5.569 | ng/ul# | 84  |
| 88) Benzo(k)fluoranthene   | 23.02 | 252 | 132138m | 1.694 | ng/ul  |     |
| 90) Benzo(a)pyrene         | 23.58 | 252 | 302118  | 3.776 | ng/ul# | 83  |
| 91) Indeno(1,2,3-cd)pyrene | 25.99 | 276 | 230491  | 2.455 | ng/ul# | 88  |
| 93) Benzo(g,h,i)perylene   | 26.71 | 276 | 158586  | 2.005 | ng/ul# | 78  |

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

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