

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\  
 Data File : BN005382.D  
 Acq On : 30 Apr 2019 21:23  
 Operator : JU/SJ  
 Sample : K2481-07ME  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 GAHW2ME

Manual Integrations  
 APPROVED

Sohil  
 5/1/2019 12:05:47 PM

Quant Time: May 01 06:58:13 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 25 03:19:34 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.59  | 152  | 336064   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.35 | 136  | 1493935  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.23 | 164  | 974197   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 16.98 | 188  | 2277281  | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.20 | 240  | 2219497  | 20.00 | ng/ul | -0.02    |
| 85) Perylene-d12          | 23.40 | 264  | 2197360  | 20.00 | ng/ul | -0.04    |

## System Monitoring Compounds

|                                |       |     |         |       |       |       |
|--------------------------------|-------|-----|---------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.22  | 96  | 34979   | 5.03  | ng/uL | 0.00  |
| 5) Phenol-d5                   | 6.78  | 99  | 739112  | 28.00 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.94  | 67  | 443502  | 26.94 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.13  | 132 | 601715  | 29.79 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.29  | 113 | 592640  | 28.72 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 8.73  | 128 | 325635  | 35.41 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 9.45  | 143 | 373319  | 40.05 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 9.98  | 165 | 655297  | 30.67 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 10.49 | 131 | 779560  | 35.38 | ng/ul | 0.00  |
| 43) Dimethylphthalate-d6       | 13.65 | 166 | 2151369 | 30.42 | ng/ul | 0.00  |
| 46) Acenaphthylene-d8          | 13.92 | 160 | 2615004 | 30.11 | ng/ul | 0.00  |
| 51) 4-Nitrophenol-d4           | 14.44 | 143 | 353614  | 31.67 | ng/ul | 0.00  |
| 57) Fluorene-d10               | 15.23 | 176 | 1866414 | 31.45 | ng/ul | 0.00  |
| 62) 4,6-Dinitro-2-methylphenol | 15.37 | 200 | 352800  | 34.45 | ng/ul | 0.01  |
| 70) Anthracene-d10             | 17.09 | 188 | 2878182 | 30.15 | ng/ul | 0.00  |
| 78) Pyrene-d10                 | 19.39 | 212 | 3428180 | 30.19 | ng/ul | 0.00  |
| 89) Benzo(a)pyrene-d12         | 23.26 | 264 | 3062961 | 31.90 | ng/ul | -0.04 |

## Target Compounds

|                            |       |     |           |         | Ovalue |    |
|----------------------------|-------|-----|-----------|---------|--------|----|
| 28) Naphthalene            | 10.40 | 128 | 1203966   | 17.336  | ng/ul  | 99 |
| 34) 2-Methylnaphthalene    | 12.03 | 142 | 11426403  | 224.697 | ng/ul  | 98 |
| 40) 1,1'-Biphenyl          | 13.05 | 154 | 1329794   | 18.320  | ng/ul  | 96 |
| 44) Dimethylphthalate      | 13.69 | 163 | 458795    | 6.511   | ng/ul  | 99 |
| 47) Acenaphthylene         | 13.95 | 152 | 7309306   | 85.630  | ng/ul# | 96 |
| 49) Acenaphthene           | 14.29 | 153 | 800173    | 13.731  | ng/ul  | 99 |
| 53) Dibenzofuran           | 14.63 | 168 | 997977    | 11.895  | ng/ul  | 95 |
| 58) Fluorene               | 15.29 | 166 | 5217946   | 79.701  | ng/ul  | 97 |
| 69) Phenanthrene           | 17.03 | 178 | 19591009m | 176.643 | ng/ul  |    |
| 71) Anthracene             | 17.12 | 178 | 2630382   | 23.282  | ng/ul  | 99 |
| 76) Fluoranthene           | 19.06 | 202 | 7355614   | 59.793  | ng/ul  | 98 |
| 79) Pyrene                 | 19.43 | 202 | 11133843  | 77.910  | ng/ul  | 98 |
| 82) Benzo(a)anthracene     | 21.19 | 228 | 4302099m  | 31.158  | ng/ul  |    |
| 84) Chrysene               | 21.25 | 228 | 4058759   | 31.673  | ng/ul  | 99 |
| 87) Benzo(b)fluoranthene   | 22.75 | 252 | 2139862   | 17.324  | ng/ul  | 99 |
| 88) Benzo(k)fluoranthene   | 22.79 | 252 | 547762m   | 4.643   | ng/ul  |    |
| 90) Benzo(a)pyrene         | 23.30 | 252 | 1332362   | 11.476  | ng/ul  | 97 |
| 91) Indeno(1,2,3-cd)pyrene | 25.57 | 276 | 654165    | 5.096   | ng/ul  | 95 |
| 92) Dibenzo(a,h)anthracene | 25.58 | 278 | 255889    | 2.349   | ng/ul  | 96 |
| 93) Benzo(g,h,i)perylene   | 26.23 | 276 | 449408    | 4.234   | ng/ul  | 95 |

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN043019\  
Data File : BN005382.D  
Acq On : 30 Apr 2019 21:23  
Operator : JU/SJ  
Sample : K2481-07ME  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
GAHW2ME

Manual Integrations  
APPROVED

Sohil  
5/1/2019 12:05:47 PM

Quant Time: May 01 06:58:13 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN041919MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Apr 25 03:19:34 2019  
Response via : Initial Calibration

