

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091219\  
 Data File : BP000217.D  
 Acq On : 12 Sep 2019 14:51  
 Operator : HP/JU  
 Sample : K4634-15DL 5X  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleID :  
 F2D33DL

Manual Integrations  
 APPROVED

mohammad  
 9/16/2019 9:09:11 AM

Quant Time: Sep 13 13:29:52 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP090519MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 10 02:43:18 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.89  | 152  | 476254   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 8.18  | 136  | 1780639  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 9.95  | 164  | 942400   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 11.46 | 188  | 1658945  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 14.17 | 240  | 1325504  | 20.00 | ng/ul | 0.02     |
| 86) Perylene-d12          | 15.75 | 264  | 1402222  | 20.00 | ng/ul | 0.06     |

## System Monitoring Compounds

|                                |       |     |        |      |       |      |
|--------------------------------|-------|-----|--------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 2.62  | 96  | 5952m  | 0.44 | ng/uL | 0.02 |
| 5) Phenol-d5                   | 6.51  | 99  | 104254 | 2.57 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.57  | 67  | 56711  | 2.67 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 6.66  | 132 | 89333  | 2.72 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 7.26  | 113 | 77939  | 2.56 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 7.45  | 128 | 37257  | 2.76 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 7.78  | 143 | 36726  | 2.79 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 8.02  | 165 | 74275  | 2.66 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 8.23  | 131 | 47709  | 1.38 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 9.63  | 166 | 216339 | 3.11 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 9.79  | 160 | 259959 | 3.02 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 10.04 | 143 | 21653  | 1.74 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 10.47 | 176 | 182100 | 3.08 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 10.54 | 200 | 5210   | 0.68 | ng/ul | 0.01 |
| 71) Anthracene-d10             | 11.51 | 188 | 259454 | 3.23 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 12.90 | 212 | 243802 | 3.09 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 15.64 | 264 | 214071 | 2.95 | ng/ul | 0.05 |

## Target Compounds

|                            |       |     |         |        | Ovalue |    |
|----------------------------|-------|-----|---------|--------|--------|----|
| 70) Phenanthrene           | 11.48 | 178 | 535388  | 5.457  | ng/ul  | 99 |
| 72) Anthracene             | 11.53 | 178 | 113347  | 1.174  | ng/ul  | 98 |
| 75) Carbazole              | 11.69 | 167 | 90182   | 1.093  | ng/ul  | 97 |
| 77) Fluoranthene           | 12.69 | 202 | 1935646 | 19.562 | ng/ul  | 99 |
| 80) Pyrene                 | 12.93 | 202 | 1896637 | 18.966 | ng/ul  | 95 |
| 83) Benzo(a)anthracene     | 14.15 | 228 | 1162184 | 12.112 | ng/ul  | 97 |
| 85) Chrysene               | 14.19 | 228 | 1217285 | 13.825 | ng/ul  | 98 |
| 88) Benzo(b)fluoranthene   | 15.27 | 252 | 988603m | 10.637 | ng/ul  |    |
| 89) Benzo(k)fluoranthene   | 15.28 | 252 | 895499m | 10.222 | ng/ul  |    |
| 91) Benzo(a)pyrene         | 15.67 | 252 | 854327  | 9.798  | ng/ul  | 99 |
| 92) Indeno(1,2,3-cd)pyrene | 17.34 | 276 | 453599m | 4.522  | ng/ul  |    |
| 93) Dibenzo(a,h)anthracene | 17.34 | 278 | 139704  | 1.681  | ng/ul# | 90 |
| 94) Benzo(g,h,i)perylene   | 17.82 | 276 | 413064  | 4.939  | ng/ul  | 97 |

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

