

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP101219\  
 Data File : BP000669.D  
 Acq On : 12 Oct 2019 15:55  
 Operator : JU  
 Sample : PB123660BL  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK60

Quant Time: Oct 12 22:52:32 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 09 17:20:45 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.75  | 152  | 174586   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 8.03  | 136  | 662438   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 9.80  | 164  | 370670   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 11.29 | 188  | 687163   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 13.97 | 240  | 619993   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 15.45 | 264  | 691443   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |       |
|--------------------------------|-------|-----|---------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 2.36  | 96  | 31266   | 7.43  | ng/uL | -0.01 |
| 5) Phenol-d5                   | 6.38  | 99  | 505796  | 37.56 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.43  | 67  | 258701  | 38.42 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 6.52  | 132 | 422369  | 37.23 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 7.13  | 113 | 394898  | 38.38 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 7.30  | 128 | 194945  | 38.71 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 7.63  | 143 | 210786  | 39.31 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 7.88  | 165 | 368796  | 35.94 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 8.09  | 131 | 518771  | 45.45 | ng/ul | 0.00  |
| 44) Dimethylphthalate-d6       | 9.49  | 166 | 1005086 | 38.73 | ng/ul | 0.00  |
| 47) Acenaphthylene-d8          | 9.64  | 160 | 1206316 | 37.79 | ng/ul | 0.00  |
| 52) 4-Nitrophenol-d4           | 9.90  | 143 | 145188  | 36.33 | ng/ul | 0.00  |
| 58) Fluorene-d10               | 10.32 | 176 | 839822  | 37.83 | ng/ul | 0.00  |
| 63) 4,6-Dinitro-2-methylphenol | 10.39 | 200 | 88735   | 27.12 | ng/ul | 0.00  |
| 71) Anthracene-d10             | 11.35 | 188 | 1259319 | 39.56 | ng/ul | 0.00  |
| 79) Pyrene-d10                 | 12.73 | 212 | 1319773 | 37.64 | ng/ul | 0.00  |
| 90) Benzo(a)pyrene-d12         | 15.36 | 264 | 1353687 | 39.66 | ng/ul | 0.00  |

Target Compounds Ovalue

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

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