

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP101719\  
 Data File : BP000843.D  
 Acq On : 17 Oct 2019 15:59  
 Operator : JU  
 Sample : PB124015BL  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK15

Quant Time: Oct 17 16:31:58 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 17 09:27:24 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.75  | 152  | 177675   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 8.03  | 136  | 677587   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 9.80  | 164  | 384140   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 11.29 | 188  | 716396   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 13.97 | 240  | 616634   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 15.45 | 264  | 693094   | 20.00 | ng/ul | 0.00     |

|                                |       |     |         |       |       |       |
|--------------------------------|-------|-----|---------|-------|-------|-------|
| System Monitoring Compounds    |       |     |         |       |       |       |
| 3) 1,4-Dioxane-d8              | 2.36  | 96  | 28352   | 6.62  | ng/uL | -0.04 |
| 5) Phenol-d5                   | 6.39  | 99  | 428132  | 31.24 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.43  | 67  | 214299  | 31.28 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 6.52  | 132 | 375211  | 32.49 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 7.13  | 113 | 342965  | 32.75 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 7.31  | 128 | 170045  | 33.01 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 7.63  | 143 | 187171  | 34.13 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 7.88  | 165 | 330675  | 31.50 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 8.09  | 131 | 461586  | 39.54 | ng/ul | 0.00  |
| 44) Dimethylphthalate-d6       | 9.49  | 166 | 887386  | 33.00 | ng/ul | 0.00  |
| 47) Acenaphthylene-d8          | 9.64  | 160 | 1062207 | 32.11 | ng/ul | 0.00  |
| 52) 4-Nitrophenol-d4           | 9.92  | 143 | 116803  | 28.20 | ng/ul | 0.00  |
| 58) Fluorene-d10               | 10.32 | 176 | 747389  | 32.48 | ng/ul | 0.00  |
| 63) 4,6-Dinitro-2-methylphenol | 10.39 | 200 | 75267   | 22.06 | ng/ul | 0.00  |
| 71) Anthracene-d10             | 11.35 | 188 | 1078976 | 32.52 | ng/ul | 0.00  |
| 79) Pyrene-d10                 | 12.74 | 212 | 1119979 | 32.12 | ng/ul | 0.00  |
| 90) Benzo(a)pyrene-d12         | 15.36 | 264 | 1119819 | 32.73 | ng/ul | 0.00  |

Target Compounds Ovalue

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

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