

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042421\  
 Data File : BN014436.D  
 Acq On : 25 Apr 2021 11:51  
 Operator : CG/JU  
 Sample : PB135972BL  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK972

Quant Time: Apr 26 05:00:47 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN041521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 23 12:06:34 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.68  | 152  | 98801    | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.46 | 136  | 436006   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.32 | 164  | 241678   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.06 | 188  | 501024   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.26 | 240  | 460334   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.48 | 264  | 470067   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.21  | 96  | 15848   | 6.11  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.60  | 84  | 204309  | 29.71 | ng/ul | 0.00 |
| 7) Phenol-d5                   | 6.85  | 99  | 321376  | 38.91 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.02  | 67  | 206927  | 41.84 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.22  | 132 | 281340  | 43.30 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.38  | 113 | 273767  | 43.73 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 8.83  | 128 | 155413  | 49.33 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.55  | 143 | 127401  | 41.29 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.08 | 165 | 268685  | 40.42 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.59 | 131 | 402672  | 40.84 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 13.73 | 166 | 887979  | 50.72 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 14.01 | 160 | 1218471 | 57.85 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.52 | 143 | 104667  | 30.82 | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.32 | 176 | 809545  | 52.31 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.43 | 200 | 60759   | 22.66 | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.16 | 188 | 1241165 | 53.08 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.46 | 212 | 1313984 | 58.80 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.34 | 264 | 1276324 | 51.18 | ng/ul | 0.00 |

Target Compounds Ovalue

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

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