

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041521\  
 Data File : BN014295.D  
 Acq On : 15 Apr 2021 22:12  
 Operator : CG/JU  
 Sample : PB135757BL  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK757

Quant Time: Apr 16 02:31:45 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN041521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 15 15:18:04 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.70  | 152  | 123640   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.47 | 136  | 499216   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.33 | 164  | 293357   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.08 | 188  | 588271   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.27 | 240  | 598775   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.50 | 264  | 654863   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.22  | 96  | 21407   | 6.59  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.62  | 84  | 279680  | 32.50 | ng/ul | 0.00 |
| 7) Phenol-d5                   | 6.86  | 99  | 350528  | 33.91 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.03  | 67  | 214382  | 34.64 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.23  | 132 | 295885  | 36.39 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.39  | 113 | 267165  | 34.10 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 8.85  | 128 | 124120  | 34.41 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.56  | 143 | 125060  | 35.40 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.09 | 165 | 256738  | 33.73 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.61 | 131 | 374613  | 33.18 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 13.75 | 166 | 774934  | 36.46 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 14.02 | 160 | 970523  | 37.96 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.52 | 143 | 113923  | 27.64 | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.33 | 176 | 673427  | 35.85 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.44 | 200 | 77699   | 24.68 | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.17 | 188 | 1054550 | 38.41 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.47 | 212 | 1217596 | 41.89 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.36 | 264 | 1409045 | 40.56 | ng/ul | 0.00 |

Target Compounds Ovalue

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

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