

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN031022\  
 Data File : BN018897.D  
 Acq On : 10 Mar 2022 10:41  
 Operator : CG/JU  
 Sample : PB142853BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK853

Quant Time: Mar 10 11:17:13 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN030822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 08 14:15:20 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.899  | 152  | 164082   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.704 | 136  | 769934   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.534 | 164  | 527253   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.275 | 188  | 1153079  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.445 | 240  | 1282749  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 23.839 | 264  | 1355851  | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.269  | 96  | 23121   | 5.636  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.687  | 84  | 339614  | 27.054 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.052  | 99  | 431396  | 26.930 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.222  | 67  | 279989  | 29.018 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.428  | 132 | 331794  | 28.901 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.604  | 113 | 349800  | 26.854 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 9.057  | 128 | 167580  | 28.336 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.787  | 143 | 183786  | 28.284 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.322 | 165 | 332433  | 26.746 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.834 | 131 | 507933  | 27.139 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.940 | 166 | 1187278 | 29.742 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.228 | 160 | 1437788 | 28.935 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.710 | 143 | 195369  | 24.166 | ng/u1 | -0.01 |
| 60) Fluorene-d10              | 15.528 | 176 | 1025108 | 29.542 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.634 | 200 | 167237  | 23.403 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.375 | 188 | 1679665 | 30.455 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.663 | 212 | 2109413 | 30.387 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.686 | 264 | 2212242 | 31.307 | ng/u1 | -0.01 |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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