

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100820\  
 Data File : BN012107.D  
 Acq On : 08 Oct 2020 17:55  
 Operator : CG/JU  
 Sample : L4253-10  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SVOC-GCP2-BLANK

Quant Time: Oct 08 18:23:53 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN100220.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 08 15:35:30 2020  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|--------|------|----------|-------|-------|----------|
| -----                     |        |      |          |       |       |          |
| Internal Standards        |        |      |          |       |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.646  | 152  | 213231   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.422 | 136  | 829565   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.287 | 164  | 515579   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.034 | 188  | 1094923  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.233 | 240  | 1122456  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.433 | 264  | 1158387  | 20.00 | ng/ul | 0.00     |

|                               |       |     |    |      |       |  |
|-------------------------------|-------|-----|----|------|-------|--|
| System Monitoring Compounds   |       |     |    |      |       |  |
| 3) 1,4-Dioxane-d8             | 0.000 | 96  | 0d | 0.00 | ng/uL |  |
| 5) Phenol-d5                  | 0.000 | 99  | 0  | 0.00 | ng/ul |  |
| 7) Bis-(2-Chloroethyl)eth...  | 0.000 | 67  | 0d | 0.00 | ng/ul |  |
| 9) 2-Chlorophenol-d4          | 0.000 | 132 | 0  | 0.00 | ng/ul |  |
| 13) 4-Methylphenol-d8         | 0.000 | 113 | 0  | 0.00 | ng/ul |  |
| 19) Nitrobenzene-d5           | 0.000 | 128 | 0  | 0.00 | ng/ul |  |
| 22) 2-Nitrophenol-d4          | 0.000 | 143 | 0  | 0.00 | ng/ul |  |
| 26) 2,4-Dichlorophenol-d3     | 0.000 | 165 | 0  | 0.00 | ng/ul |  |
| 29) 4-Chloroaniline-d4        | 0.000 | 131 | 0d | 0.00 | ng/ul |  |
| 44) Dimethylphthalate-d6      | 0.000 | 166 | 0  | 0.00 | ng/ul |  |
| 47) Acenaphthylene-d8         | 0.000 | 160 | 0  | 0.00 | ng/ul |  |
| 52) 4-Nitrophenol-d4          | 0.000 | 143 | 0  | 0.00 | ng/ul |  |
| 58) Fluorene-d10              | 0.000 | 176 | 0  | 0.00 | ng/ul |  |
| 63) 4,6-Dinitro-2-methylph... | 0.000 | 200 | 0  | 0.00 | ng/ul |  |
| 71) Anthracene-d10            | 0.000 | 188 | 0d | 0.00 | ng/ul |  |
| 79) Pyrene-d10                | 0.000 | 212 | 0d | 0.00 | ng/ul |  |
| 90) Benzo(a)pyrene-d12        | 0.000 | 264 | 0d | 0.00 | ng/ul |  |

Target Compounds Qvalue

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

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